# FRENCH LIMITED SITE CROSBY, TEXAS

# DNAPL INVESTIGATION PRELIMINARY REPORT

Submitted to:

FLTG, Inc. Crosby, Texas

2002953



Submitted by:

Applied Hydrology Associates, Inc. Denver, Colorado

July 1992

01502432

#### TABLE OF CONTENTS

			Page
1.0	Introd	luction	1
	1.1	Background	1
	1.2	Conceptual Model of DNAPL Migration	2
2.0	Aprıl	1992 Investigation	3
	2.1	S1-16 Area	3
	2.2	INT-11 Area	4
	2.3	Area Inside Floodwall	6
3.0	Summ	nary and Conclusions	7

#### **Figures**

- 1 DNAPL Investigation Location Map
- 2 Location Map, DNAPL Investigation, S1-16 Area
- 3 Cross Sections A-A' & B-B', DNAPL Investigation, S1-16 Area
- 4 Structure on the C1 Interval, DNAPL Investigation, S1-16 Area
- 5 C1 Clay Isopach (Ratio % > 3), DNAPL Investigation, S1-16 Area
- 6 DNAPL Occurrence in the S1 interval, DNAPL Investigation, S1-16 Area
- 7 Location Map, DNAPL Investigation, INT-11 Area
- 8 Cross Section C-C', DNAPL Investigation, INT-11 Area
- 9 C1 Clay Isopach (Ratio % > 3), DNAPL Investigation, INT-11 Area
- 10 Structure on the C1 Interval, DNAPL Investigation, INT-11 Area
- 11 Structure on the C2 Interval, DNAPL Investigation, INT-11 Area
- 12 DNAPL Occurrence in the INT interval, DNAPL Investigation, INT-11 Area

#### **Tables**

- 1 Summary of Soil Sample Analyses, S1-16 and INT-11 Areas
- 2 Sample Descriptions, S1-16 and INT-11 Areas
- 3 Volatile Organic Analyses, Well INT-64, INT-11 Area
- 4 Analysis of Soil Samples from INT-11B and INT-11C

#### **Appendices**

- 1 Analytical Reports S1-16 DNAPL samples
- 2 Analytical Reports INT-11 DNAPL sample
- 3 Cone Penetrometer Logs S1-16 Area
- 4 Cone Penetrometer Logs INT-11 Area
- 5 Boring Log and Well Construction Details for REI-6-1
- 6 Analytical Reports INT-11B and INT-11C Soil samples

#### 1.0 INTRODUCTION

This report describes the detailed results of an initial investigation conducted in the S1-16 and INT-11 well areas in April 1992 to delineate the extent of free-phase organic liquid. Major results of this investigation were first presented in the April 1992 Monthly Progress Report.

#### 1.1 Background

The geology of the French Limited site is divided into three zones<sup>1</sup>: the lower silty sand zone, the middle clayey zone, and the upper alluvial zone. The lower silty sand zone is a confined aquifer, the top of which is between 109 and 137 feet below ground surface (BGS). It is separated from the upper alluvial zone by the middle clayey zone, a 60- to 90-foot-thick aquitard, referred to as the C2 clay. The upper alluvial zone is up to 55 feet thick. It is subdivided into a lower water-bearing zone of interbedded sands, silts, and clays (the INT unit), an upper sandy water-bearing zone (the S1 unit), and a clay aquitard (the C1 unit) that separates the INT and S1 units. Contaminated groundwater at the site is limited to the S1 and INT units, and is being remediated by a combination of source cleanup (lagoon bioremediation), physical and hydraulic source migration controls (a sheet-pile migration control wall and an inward-gradient groundwater extraction system), groundwater extraction and surface treatment using a fixed-film bioreactor, and in-situ bioremediation using a system of injection wells.

Dense non-aqueous phase liquids (DNAPLs) were first detected at the S1-16 and INT-11 production wells in January, 1992, shortly after initiation of the production well system in early January 1992. Well S1-16 is located inside the floodwall at the southeast end of the lagoon: well INT-11 is located outside the floodwall, east of center of the south side of the lagoon (see Figure 1).

A sample of DNAPL was collected from well S1-16 in January 1992 and analyzed for volatile organic compounds (VOCs), semi-volatile organic compounds (SVOCs), total organic carbon (TOC), total petroleum hydrocarbons (TPH), and total chlorinated hydrocarbons (TOX). The results of these analyses are presented in Appendix 1. The sample contained very high concentrations of VOCs including chloroform (7.8%), 1,2-DCA (6.5%), carbon tetrachloride (2%), 1,1-DCA (0.2%), 1,2-DCE (0.2%), and TCE (0.1%), and tentatively identified compounds (TICs) including bromochloromethane (2.5%) and 1,4-difluorobenzene (2.3%). Therefore, total VOCs exceeded 20%. The sample also contained high concentrations of SVOCs including naphthalene (0.5%).

A sample of DNAPL was collected from well INT-11 in March 1992 and analyzed for VOCs, TOC and TOX. The results of these analyses are presented in Appendix 2. The sample contained high concentrations of VOCs including chloroform (810,000 ppb), 1,2-DCA (650,000 ppb), vinyl acetate (150,000 ppb), 4-methyl-2-pentanone (27,000 ppb), 2-butanone (23,000 ppb), trans-1,2-DCE (21,000 ppb), and other VOCs detected at 20,000 ppb or less. Total VOCs detected totaled 0.18%.

The April 1992 investigation was intended to follow up these indications of DNAPLs, and was conducted in two areas: around well S1-16, and around well INT-11. The aim of the investigation was to define the extent of free DNAPL outside the floodwall. The two areas investigated are shown on Figure 1.

<sup>&</sup>lt;sup>1</sup> For more details, see Hydrogeologic Assessment Report, March, 1989, Chapters 2 and 3

# 1.2 Conceptual model of DNAPL migration

There are two mechanisms that may account for the presence of DNAPLs in borings near the French Lagoon:

- 1) Migration as a separate phase directly from the lagoon; and/or
- 2) Mobilization of organic constituents near the lagoon due to their increased solubility in the presence of organic solvents, followed by deposition further away from the lagoon due to decreasing solvent concentrations and lower solubility<sup>2</sup>.

Because DNAPLs are denser than groundwater, and can be retained by aquifer matrix material, DNAPL migration below the water table at the French Limited site is principally controlled by the following factors<sup>3</sup>:

- 1) Residual saturation of aquifer matrix, i.e., the amount of DNAPL that is retained ("residual DNAPL"):
- 2) Direction and hydraulic gradient of groundwater flow, causing unretained ("free") DNAPL to migrate downgradient;
- 3) Slope of the base surface of aquifer units, causing free DNAPL that has sunk to the base of the aquifer to migrate down-dip under gravity; and
- 4) Construction of the sheet-pile flood and migration control wall ("floodwall") in 1989, which, together with inward-gradient pumping from production wells inside the floodwall, should prevent any further migration of groundwater and DNAPL beyond the floodwall.

At the site, the present extent of free and residual DNAPL outside the floodwall likely reflects conditions existing before lagoon remediation and floodwall construction. The extent of free DNAPL within the floodwall probably has been modified by floodwall construction and inward-gradient pumping from wells within the floodwall. DNAPL present beneath the French lagoon has likely migrated downward to the top of the low permeability zone within and at the base of the S1 and INT units, down the dip slope of these low-permeability zones, and, where influenced by stronger hydraulic gradients near production wells, toward the wells pumping from inside the floodwall. However, the floodwall barrier is likely to prevent wells pumping inside the floodwall drawing free DNAPL back from outside the floodwall.

The main geological features of the site likely to affect DNAPL migration are the C1 and C2 clay units that divide and underlie the upper alluvial zone, respectively. The C1 unit separates the overlying S1 sandy unit from the underlying INT interbedded unit. The C2 unit is a thick clay that underlies the INT unit. The two clay units are probable effective barriers to DNAPL migration. However, as described below, the C1 unit thickness varies across the site and it is absent under part of the site. Subordinate geological features likely to retard downward DNAPL migration are minor clay layers within the S1 and INT units.

<sup>&</sup>lt;sup>2</sup> This mechanism is described in more detail in *Hydrogeologic Characterization Report*, March 1989, Section 4.1 1.1 Note that, under this mechanism, because high-VOC water from the lagoon is denser than ambient groundwater, it would migrate under density difference in a similar manner to DNAPL.

<sup>&</sup>lt;sup>3</sup> Discussed further in Schwille, F, 1988,. Dense Chlorinated Solvents in Porous and Fractured Media, Model Experiments, Lewis Publishers, 146 pp See pp 5-8 and figure IV b for a description of the effects of density, matrix retention, and sloping aquifer base

#### 2.0 APRIL 1992 INVESTIGATION

The DNAPL investigation in the S1-16 and INT-11 areas was conducted using a cone-penetrometer test (CPT) rig. The rig was used to perform standard CPT soundings which provide detailed vertical engineering and lithologic profiles. This equipment measures the resistance of geologic materials to the penetration of a cone tool as it is pushed into the ground hydraulically. Both the tip resistance and the sleeve or friction resistance to penetration are recorded at intervals of 0.2 feet. Lithologies are defined by the tip and friction resistances and the ratio of these values. Sand units tend to have relatively high tip resistance but low friction resistance compared to clay units. Generally, a friction to tip resistance ratio of greater than 3% indicates clay soils.

A conductivity tool was also run in conjunction with the standard cone-penetrometer tool. The conductivity tool measures the electrical conductance (EC) of the geologic unit, which is a function of both the intergranular fluid and the soil matrix. DNAPLs have very low EC relative to groundwater, particularly high-TDS groundwater. Hence, zones containing a high content of DNAPL in the pore spaces would have a correspondingly lower overall EC than the same soil containing more conductive groundwater. However, if the soil pores contain a mixture of DNAPL and high-TDS groundwater, this effect could be masked. As described in Section 1.2, it is likely that DNAPL and denser groundwater would follow similar migration paths. Another potentially complicating factor is that sandy soils have a lower EC relative to clay soils. Both types of soil are typical of the site geology. These variations in soil type result in "background" variations in EC that must be considered in interpreting CPT-EC profiles.

After running the combination logs, the resistance, ratio, and conductivity profiles were interpreted to determine the depth zones at which DNAPL occurrence was most likely. The CPT rig was then used in soil sampling mode. In this mode, the rig is used to collect narrow-diameter split-spoon soil samples from the selected depth zones, in the same area as the CPT-EC soundings. The soil samples collected are about 1 inch in diameter and 1.5 feet in length. All CPT and soil sampling holes were cemented immediately after completion using cement grout tremied to the bottom of the hole.

During the April 1992 fieldwork, all CPT and soil sampling locations were surveyed for location coordinates and ground surface elevation. Borings and CPTs from earlier investigations at the site were incorporated into the evaluation of the S1-16 and INT-11 areas.

#### 2.1 S1-16 Area

The locations of CPT soundings and soil samples in the S1-16 area are shown on Figure 2. CPT logs are presented in Appendix 3 Interpreted geologic cross sections A-A' and B-B' through the S1-16 area are shown in Figure 3. Both cross sections show the S1 sand unit, underlain by several feet of C1 clay. The relatively thick C1 Clay underlying this area is a probable competent low-permeability barrier to downward DNAPL migration.

Cross sections A-A' and B-B' show the CPT ratio percent profiles arranged in southwest to northeast transects across the investigation area, adjusted to sea level datum. Where available, conductivity curves are superimposed on ratio percent profiles. The ratio percent profiles indicate a thick, fairly uniform sand layer in the S1 interval. Descriptions of samples collected from this interval typically report a fining-upward sequence of coarse, occasionally gravelly sands at the base of the S1, grading to fine to medium sands at the top of the S1. The unit thickens from west to east across the S1-16 area.

Conductivity measurements in the S1 generally increase from top to bottom at individual CPT locations, and also from northeast to southwest in the S1 interval along both cross section traverses. Because the ratio percent profile does not indicate clay layering within the S1, the higher conductivities are likely indicative of higher TDS pore waters. Therefore, the EC pattern probably reflects the presence of high-TDS waters in the S1 unit to the south of French Lagoon and the migration of these denser fluids downward and along the base of the S1 sand.

From the interpreted CPT logs, the elevation of the top of the C1 clay (base of the S1 sand) was determined. Figure 4 shows the structure contours of this surface in the S1-16 area. The structure map was constructed based on CPT soundings performed during the April 1992 fieldwork, and CPT soundings performed during earlier investigations at the site. The top of the C1 generally dips to the east and southeast across the investigation area, away from the French Lagoon source. The elevation of the surface varies from estimated highs of -10.3 feet at C-11 and -11.8 feet at S1-16A to a low of -18.6 at S1-16D (exact elevations of the C1 surface are not available at some locations because the ground surface was not surveyed at the time of the earlier [pre-April 1992] CPT sounding, or was subsequently altered due to addition of fill, etc.). The dip of the C1 surface is approximately 0.13 ft/ft.

The C1 Clay isopach contour map (Figure 5) shows the thickness of the C1 unit beneath the S1-16 area. Figure 5 shows that the C1 clay generally exceeds 3 feet in thickness in this area, and thickens to the east and north. The C1 also thickens to the south outside the area of investigation. No data is available on the C1 thickness under French Lagoon.

Figure 6 summarizes data obtained regarding DNAPL occurrence in the S1 interval in the S1-16 area. This figure includes data obtained before and during the April 1992 investigation. These include analyses of soil samples collected from S1-17 (included in Table 1), and field observations of oily stains, residues, odors, etc. made during drilling (observations at S1-16A, S1-16C, S1-16G, and S1-16H are included in Table 2). Also shown in Figure 6 are organic vapor meter (OVM) readings made during drilling, and sites where conductivity anomalies (zones of relatively lower EC) were noted.

Inside the flood wall, CPT-EC sounding S1-16A was conducted near well S1-16 where DNAPL was detected. A conductivity anomaly close to the base of the S1 unit can be seen in Figure 3. Fluid samples retrieved from this interval using a small bailer lowered through the cone rods did not contain any DNAPL. However, during cementing of the CPT hole, a black only mixture was observed in the cement returns, indicating that DNAPL had been present in the hole. Due to the high viscosity of DNAPL, it is difficult to retrieve by bailing. A second CPT-EC sounding conducted inside the floodwall at location S1-16B, east of the lagoon, did not indicate a low-conductivity zone, and no discoloration of the cement was noted during grouting.

Outside the floodwall, cone-penetrometer profiles were conducted at approximately 25 ft spacing in the vicinity of well S1-16. EC anomalies were noted at S1-16D and S1-16E. Because the top of the C1 unit slopes to the southeast, a number of split-spoon samples were attempted in the area outside the floodwall and southeast of S1-16. Unfortunately, attempts to retrieve soil samples at the S1/C1 interface were largely unsuccessful. The sampler apparently could not retain the fairly coarse-grained material typical of the base of the S1 unit. The only successful soil retrievals were at locations S1-16G and S1-16H. The soil samples appeared to contain residual DNAPL at the bottom of the S1.

Mainly because the sampling of the lower S1 material was not successful, further work is required in this area to clearly define the extent of DNAPL occurrence. Outside the floodwall the extent of DNAPL is defined to the northeast at S1-16H, and to the southwest at S1-16F. However, it is not defined to the southeast. The northwestern extent of this DNAPL is not defined but is likely the lagoon source.

## 2.2 INT-11 Area

The locations of CPT-EC soundings and soil samples in the INT-11 area are shown on Figure 7. CPT logs are presented in Appendix 4. An interpreted east-west geologic cross section through the INT-11 area, section C-C', prepared in the same way as sections A-A' and B-B', is shown in Figure 8. It shows the S1 sand unit overlying the INT interval, separated from it by a variable thickness of C1 clay. The C1 clay thickness decreases from 6 6 feet at INT-64 to 1 foot or less at INT-10, S1-11, and INT-63, and continues to thin to the west. The C1 clay isopach contour map (Figure 9) shows the thickness of the C1 unit beneath the INT-11 area. Figure 9 shows that the C1 clay is thin (<1 foot) to the north and west of

the INT-11 area, and is absent west of INT-10. The thinning in the C1 clay appears to be due to erosion and downcutting during deposition of the S1 sand. Clearly, where the C1 clay is thin or absent it is unlikely to act as a barrier to downward DNAPL migration.

It is also possible that abandoned well REI-6-2 (see Figure 12 for location) could have acted as an additional vertical conduit for DNAPL. The construction details of REI-6-2 are presented in Appendix 5 and are shown projected onto section C-C' in Figure 8. The projected top of the water intake interval (comprising the screen and filter pack) is close to the top of the C1 clay. Although the boring log indicates 4 feet of clay at this location, adjacent CPT soundings show a much thinner clay. Therefore, it is possible that the water intake interval could connect the S1 and INT units. Because it is no longer in use, this well should be properly plugged and abandoned during future drilling work at the site.

From the interpreted CPT logs, the elevation of the top of the C1 clay (base of the S1 sand) was determined. Figure 10 shows the structure contours of this surface in the INT-11 area. The top of the C1 clay forms a low area under and south of the French Lagoon in the north central portion of the INT-11 area. This low corresponds with the area of thinning in the C1 clay described above, which is consistent with the removal of the C1 clay being due to erosion and downcutting during deposition of the S1 sand. Therefore, the top of the C1 generally dips inward toward the S1-11 and INT-11 locations, which would tend to cause DNAPL, if present, to accumulate near S1-11 and INT-11, in the area where the C1 clay is thin. The elevation of the surface varies from estimated highs of -10.8 feet at INT-64, and -13 feet between INT-10 and S1-23, to lows of -17.6 at INT-11 and -18.6 at S1-11 (exact elevations of the C1 surface are not available at some locations because the ground surface was not surveyed at the time of the earlier [pre-April 1992] CPT sounding, or was subsequently altered due to addition of fill, etc.) The inward dip of the C1 surface is approximately 0.11 ft/ft.

Where the C1 clay is thin or absent, and therefore unlikely to act as a barrier to downward migration, DNAPLs would tend to migrate downward to the top of the C2 clay (Beaumount Formation). The elevation of the top of the C2 clay (base of the INT interval) is shown on Figure 11. Figure 11 shows that the C2 clay in this area mainly dips to the east-northeast, from an estimated high of -31.5 feet at INT-63 to a low of -36.1 feet in INT-64. This is the main direction toward which DNAPL reaching the top of the C2 in this area would be expected to migrate. The dip of the C2 surface is approximately 0.04 ft/ft. To the west of the INT-11 area, the C2 surface dips at a similar gradient to the west-southwest. from a ridge between C-3 and INT-63.

Figure 12 summarizes data obtained regarding DNAPL occurrence in the INT-11 area. This figure includes data obtained before and during the April 1992 investigation. These include TOC and TPH analyses of soil samples collected from INT-10 and INT-64 (included in Table 1), and field observations of oily stains, residues, odors, etc. made during CPT soil sampling (observations at INT-11A through INT-11G are included in Table 2). VOC analyses of soil samples collected from INT-64 are presented in Table 3. Analyses of soil samples collected at INT-11B and INT-11C are included in Appendix 6 and summarized in Table 4. Also shown in Figure 12 are organic vapor meter (OVM) readings made during drilling, and sites where conductivity anomalies (zones of relatively lower EC) were noted.

A CPT-EC sounding (INT-11A) was conducted near the INT-11 well where DNAPL is known to exist. A conductivity anomaly close to the base of the INT unit at INT-11A can be seen in Figure 8. The conductivity profile indicates a zone of relatively lower conductivity close to the base of the INT unit just above the underlying C2 Clay. Fluid samples retrieved through the cone rods using a small bailer from this interval did not show any DNAPL. As indicated above, it is difficult to retrieve the higher viscosity free-phase fluids by this method. However, during cementing of the CPT hole, a black oily mixture was observed in the cement returns indicating the presence of DNAPL in this hole. Field observations made during CPT soil sampling also suggested possible free DNAPL at INT-11B through F. Soil samples obtained from the 48 to 49 ft depth interval at these locations showed evidence of DNAPL residual. The soil samples left an oily residue on the core barrel and when squeezed the soils released a watery oily fluid. However, free DNAPL in the pore spaces of these soil samples was not evident

Field observations made during drilling indicated residual DNAPL at INT-10. However, soil samples collected from INT-10 did not contain detectable TOC. VOC analyses were not performed on these samples. INT-10 may represent an area of residual DNAPL. Soil samples collected from INT-64 did not contain detectable TPH and did not show staining or free DNAPL. However, soil samples collected from INT-64 (Table 3) showed high concentrations of VOCs, particularly in the lower 10 feet, i.e., above the C2 clay. VOCs detected include methylene chloride, (5.1 ppm), 1,2-dichloroethane (4.4 ppm), acetone (3.4 ppm), carbon disulfide (2.8 ppm), and chloroform (1 ppm). These high concentrations of VOCs and their distribution with depth suggest that INT-64 may be at the fringe of DNAPL migration from the INT-11 area.

Soil samples were collected from INT-11B and INT-11C locations (Table 4). Including tentatively identified compounds and compounds identified below the detection limit, the soil samples from INT-11B contained 4.7 to 8.4 percent VOCs and 6.1 percent SVOCs; and from INT-11C, 3.2 to 3.3 percent VOCs and 2 percent SVOCs. The major VOC components are chloroform, 1,2-DCE, 1,2-DCA, carbon tetrachloride, and PCE.

The field observations and laboratory data presented support the conclusion that DNAPL from the lagoon area accumulated at the base of the S1 sand in the S1-11/INT-11 area, migrated downward through the thin or missing C1 clay (and, potentially, abandoned well REI-6-1) to the base of the INT interval, and then migrated down the dip of the C2 clay, mainly to the east-northeast. DNAPL residuals have been observed 10-15 feet to the south of INT-11 and are present over much of the area bounded by the INT-11 well, the floodgate, the sheetpile floodwall and the property fence. CPT soundings could not be completed south of the property boundary due to access difficulties. The extent of DNAPL appears to be defined to the east by its absence at S1-22, DB-19, and INT-11G, and to the southwest by its absence at DB-18. However, it is not defined to the south of INT-11

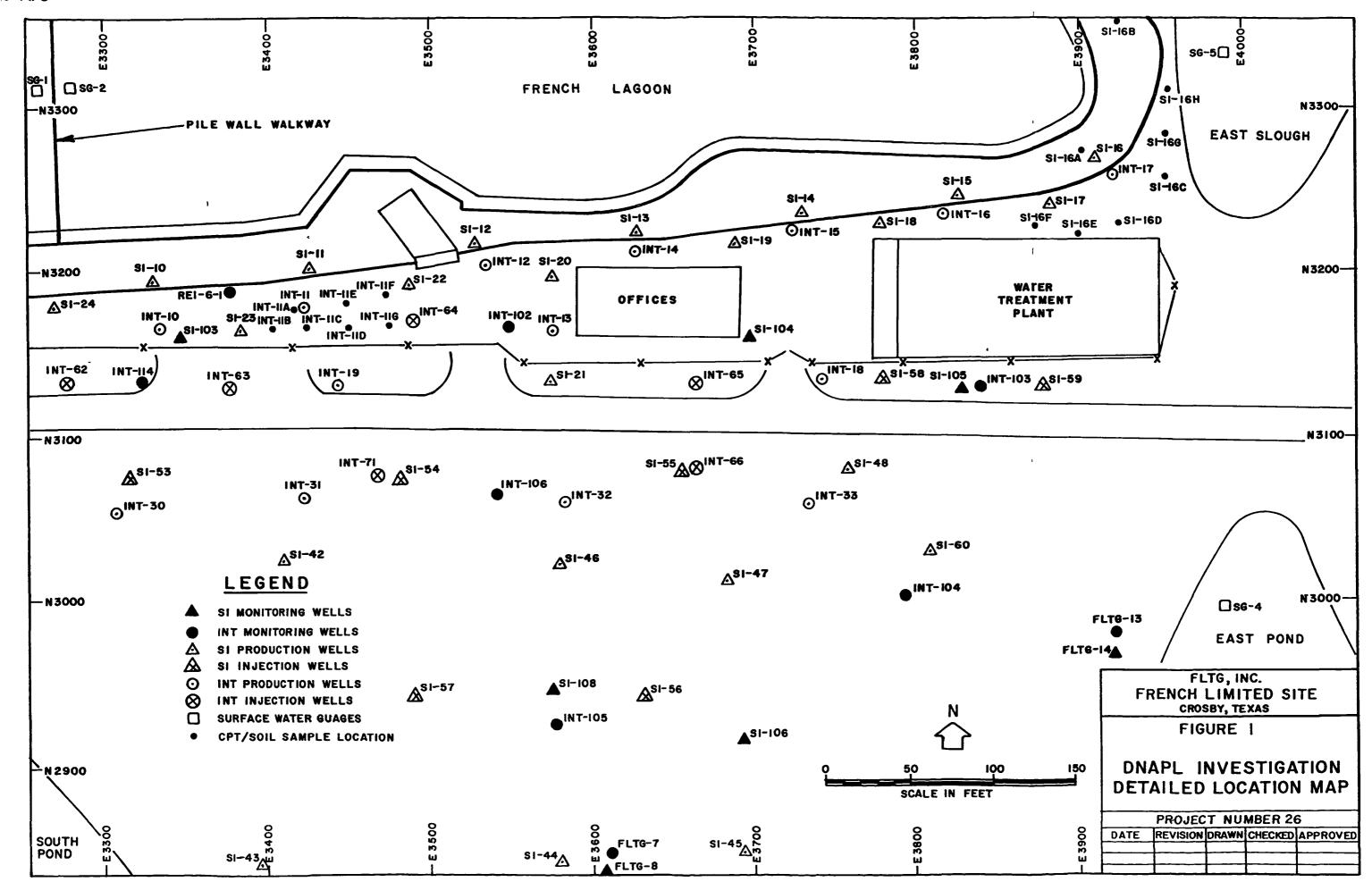
At INT-10, a less well-defined occurrence of DNAPL is indicated only by staining and odor (Figure 12). In this area, the C2 surface dips toward the west-southwest (Figure 11). DNAPLs probably entered the INT interval in the same manner as described above and followed the local structural gradient at the base of the INT interval toward INT-10 and DB-17.

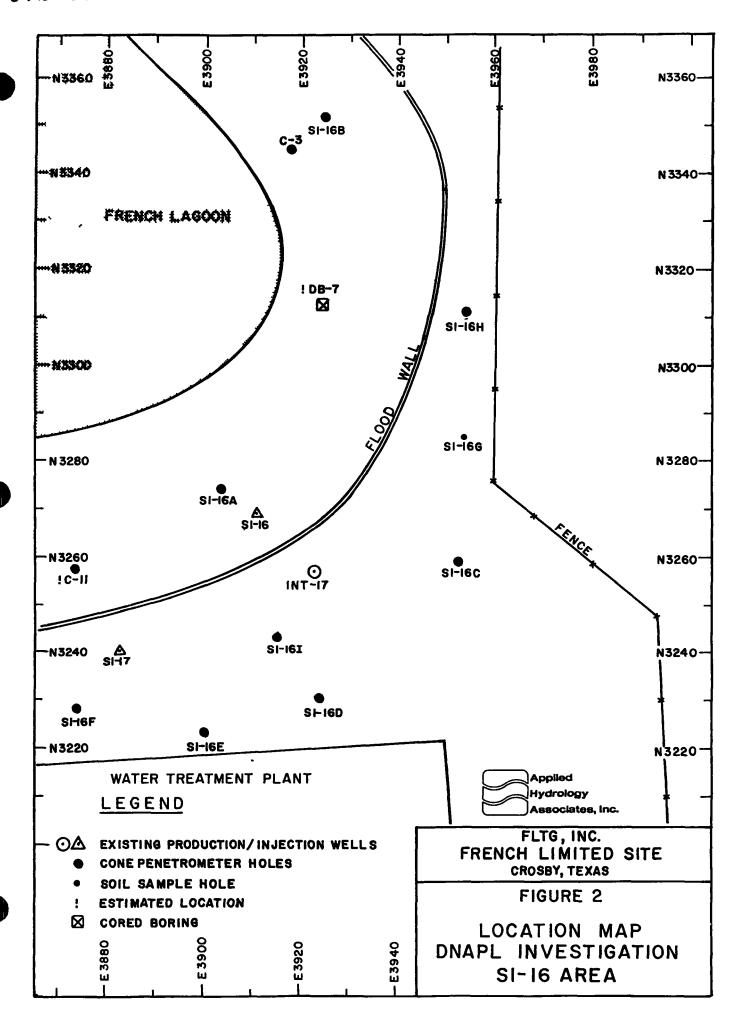
#### 2.3 Area Inside Floodwall

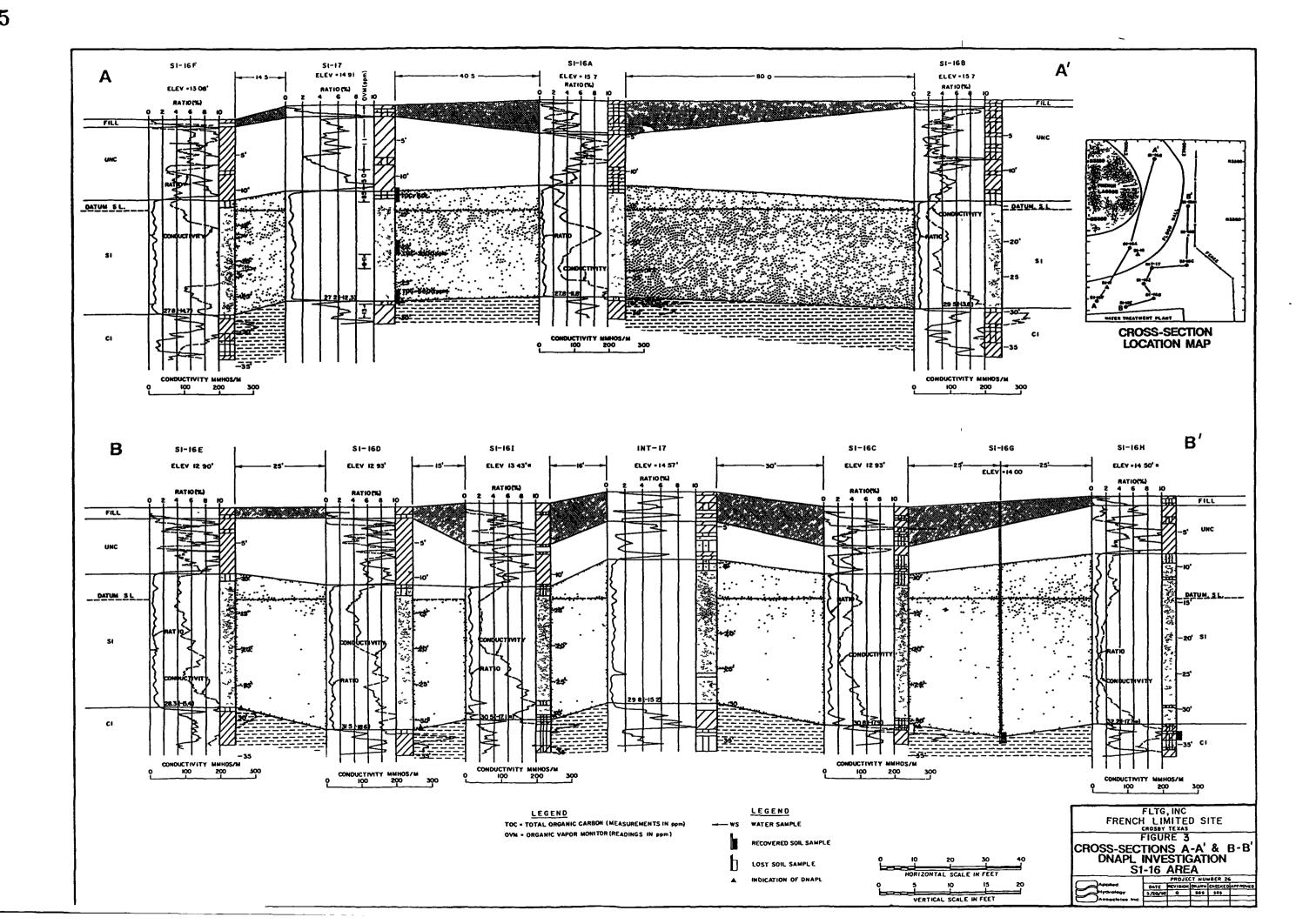
Additional observations were made during operation of the production well system after the April 1992 field investigation was complete. Since the April 1992 investigation, DNAPL has been reported at more wells inside the flood wall. In June 1992, DNAPL was observed in clogged well-head flow meters, pressure gauges, or bends in piping further down the system, at wells S1-4, S1-13, and S1-14. Samples from S1-4, S1-13, and S1-14 have been submitted for analysis for VOCs, SVOCs, TOC, and TOX. The detection of DNAPL at these wells is consistent with the conceptual model of DNAPL migration presented above, i.e., operation of the production wells inside the floodwall has pulled DNAPL previously underlying the lagoon toward those wells.

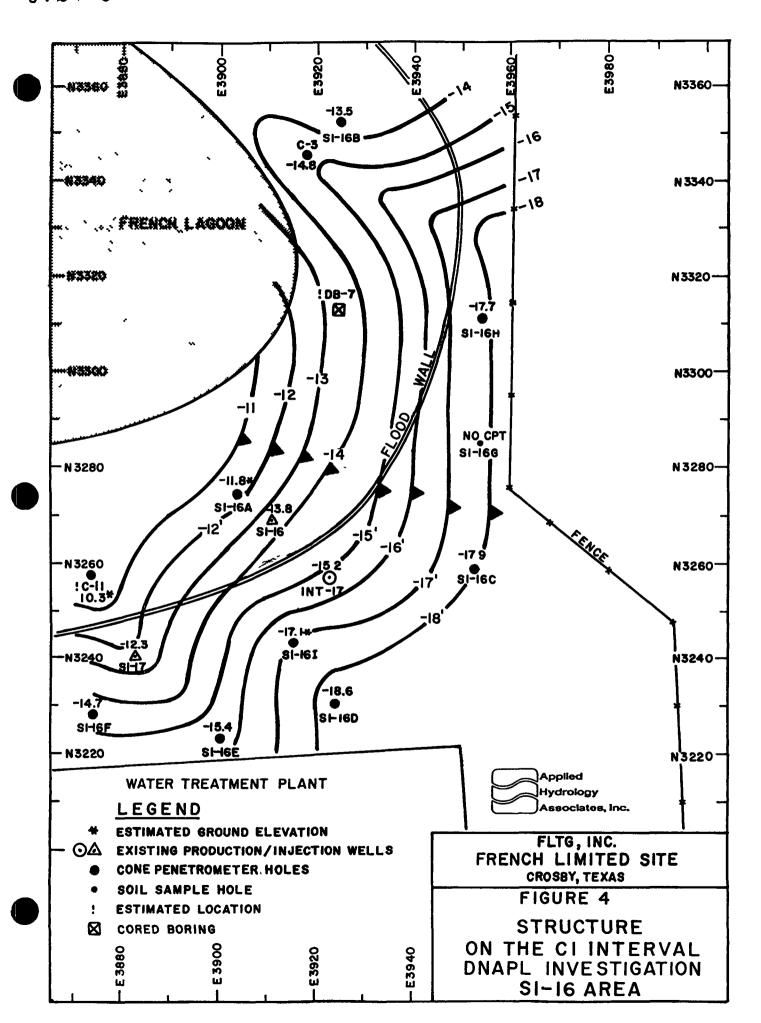
#### 3.0 SUMMARY AND CONCLUSIONS

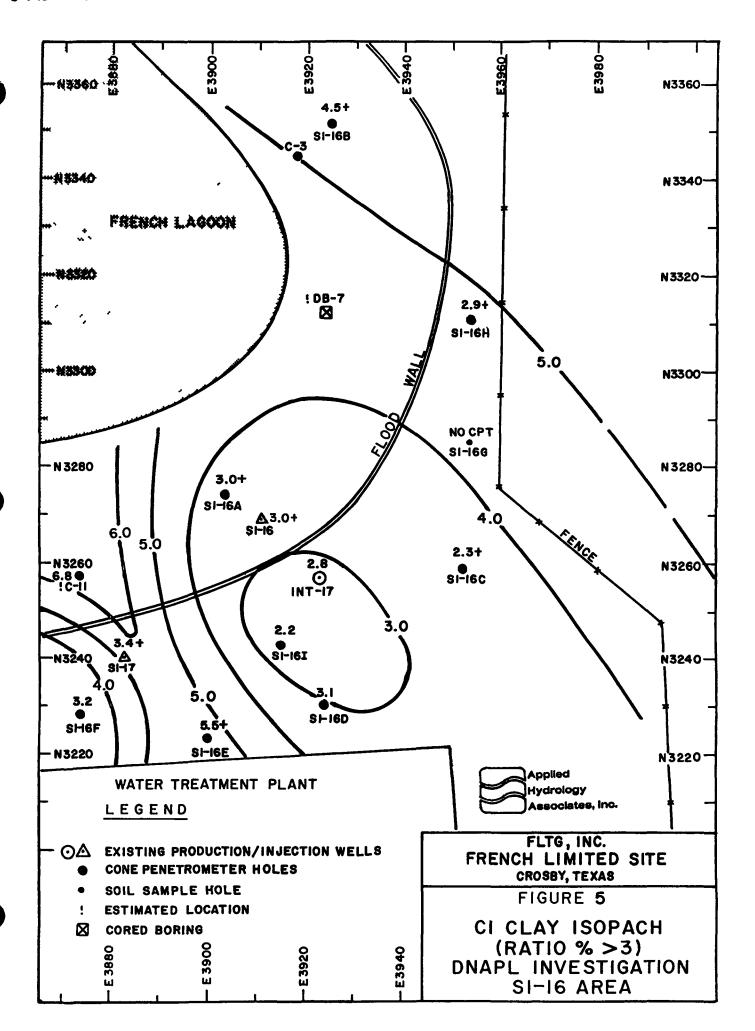
- The April, 1992 field investigations combined with data from earlier investigations, confirms the
  presence of DNAPL residuals in soils outside the floodwall in the S1-16 and INT-11 areas. DNAPL
  appears to be concentrated at the base of the S1 unit in the S1-16 area, and at the base of the INT
  unit in the INT-11 area. Similar suites of VOCs are detected in impacted soil samples, and in
  DNAPL samples.
- 2. The ocurrence and distribution of observed DNAPL is consistent with DNAPL migration controlled primarily by the structural configuration of the C1 and C2 clay units.
- 3. More detailed soil and groundwater sampling needs to be performed in the S1-16 and INT-11 areas to fully define DNAPL distribution in these areas. Proposed locations for additional borings and DNAPL detection wells are shown on Figure 1. Details of proposed soil sampling and well installation will be presented in the DNAPL Additional Investigation Work Plan.

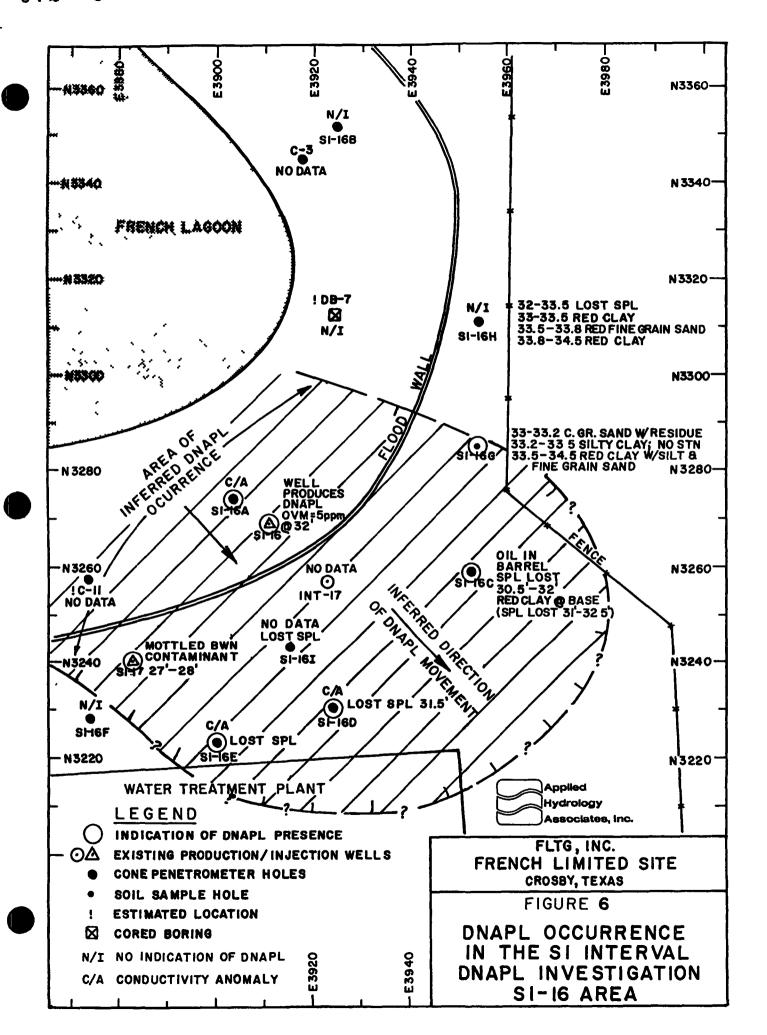


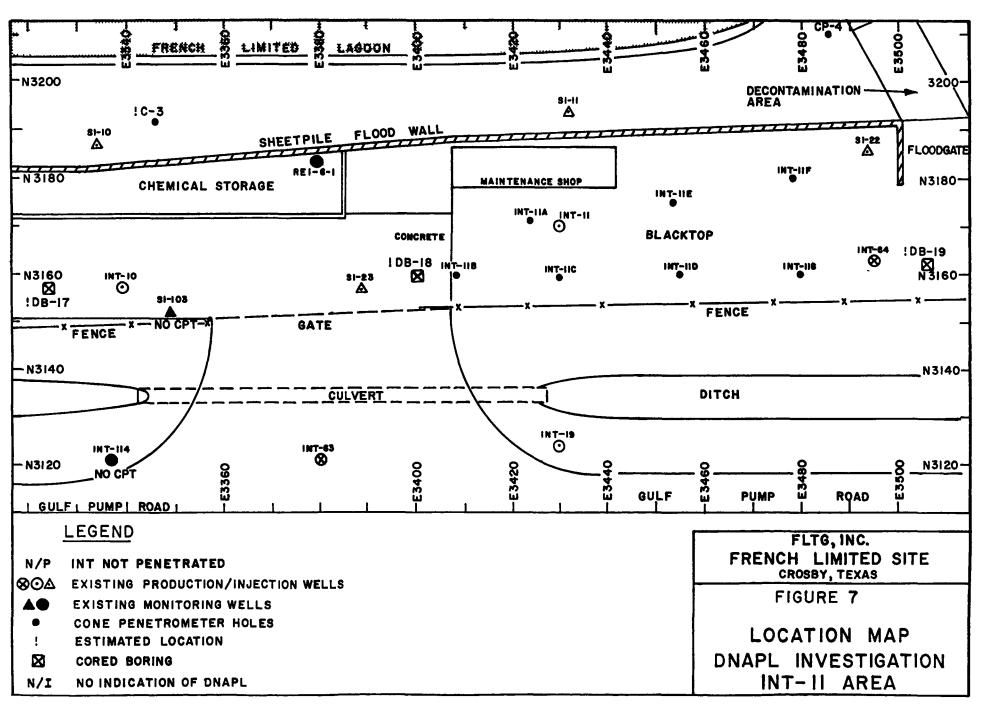


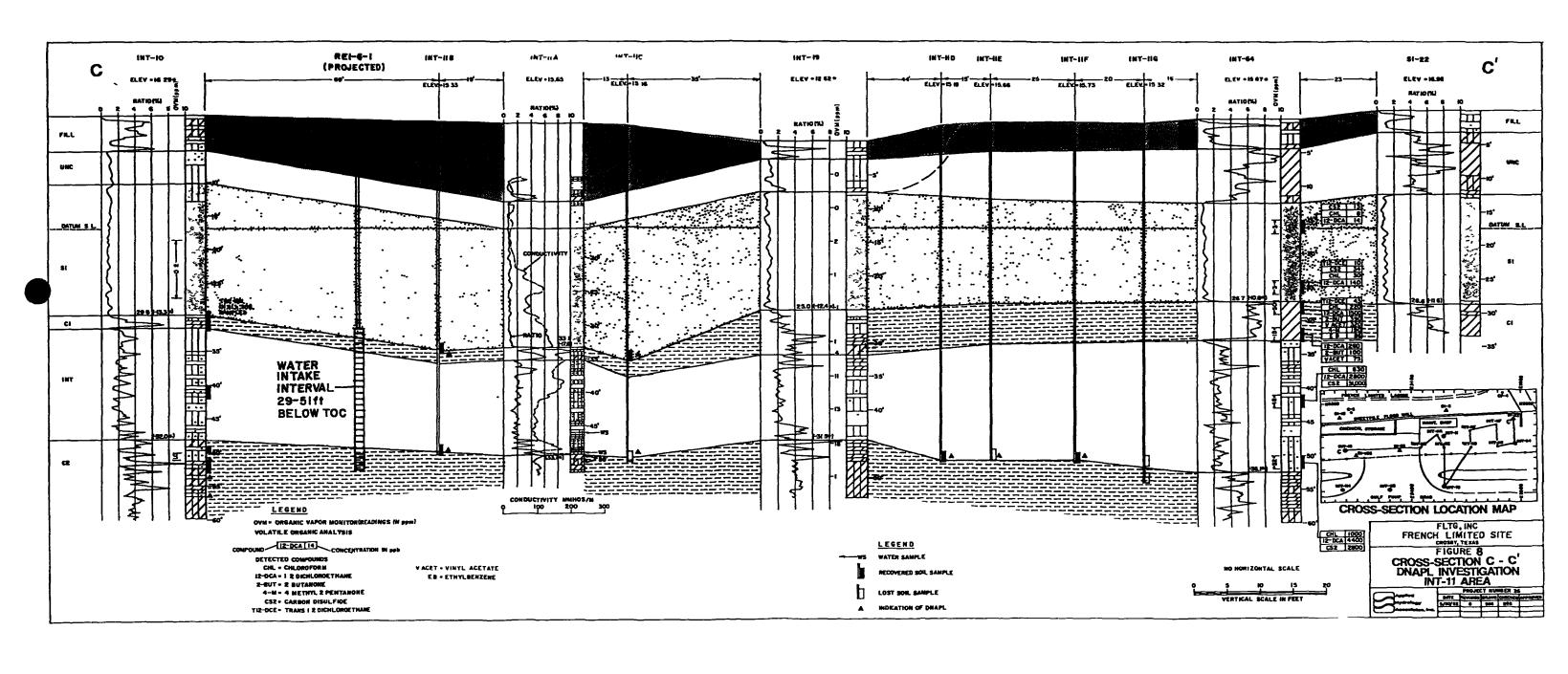


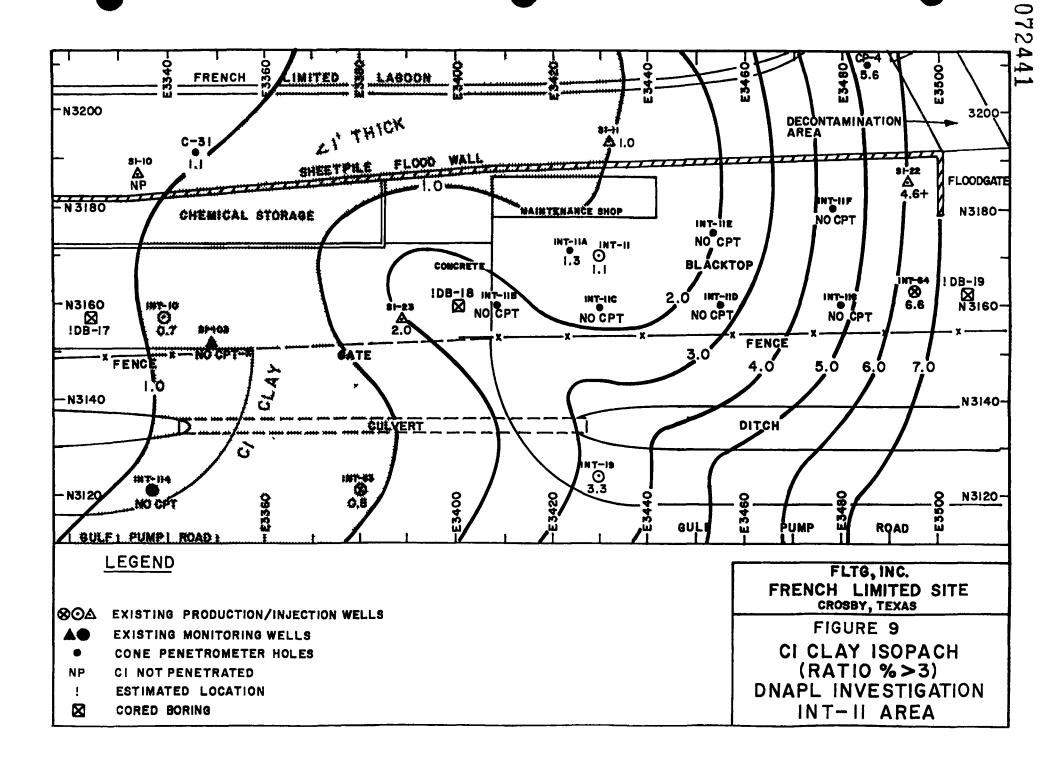


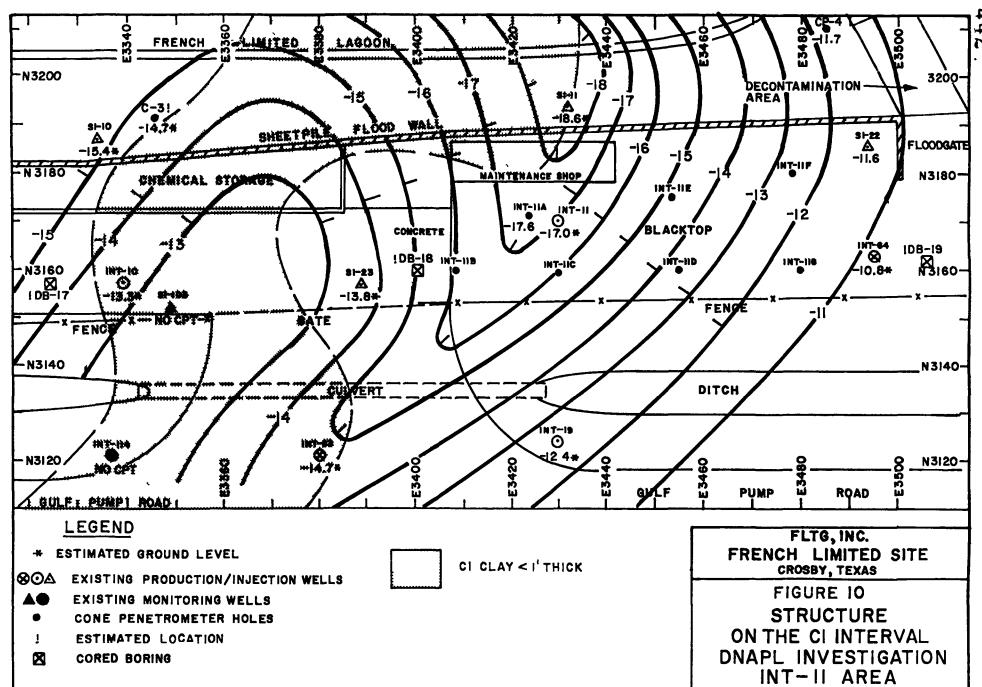


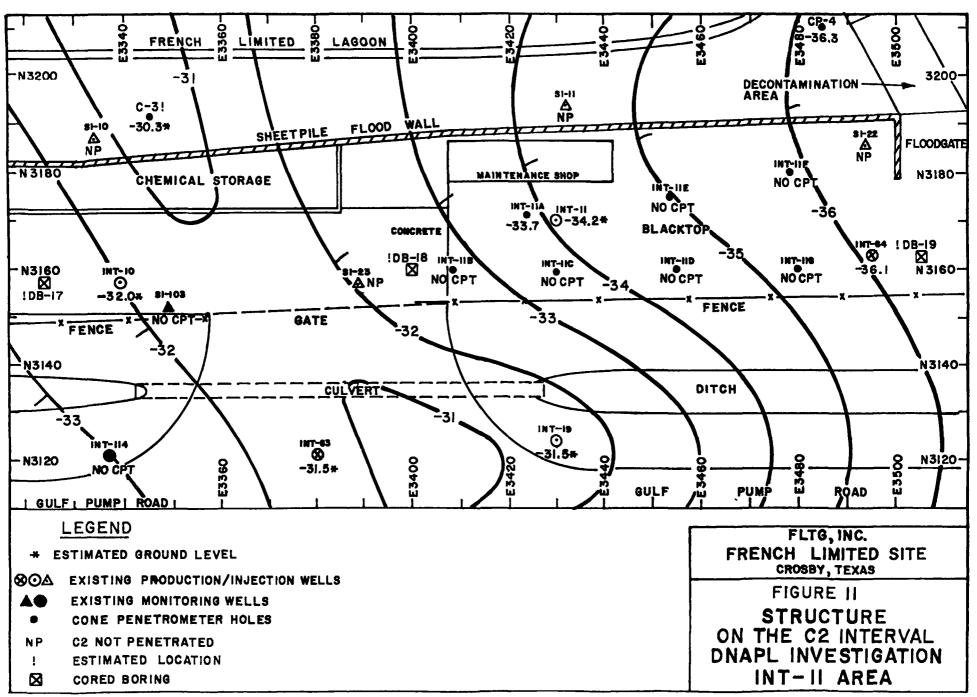












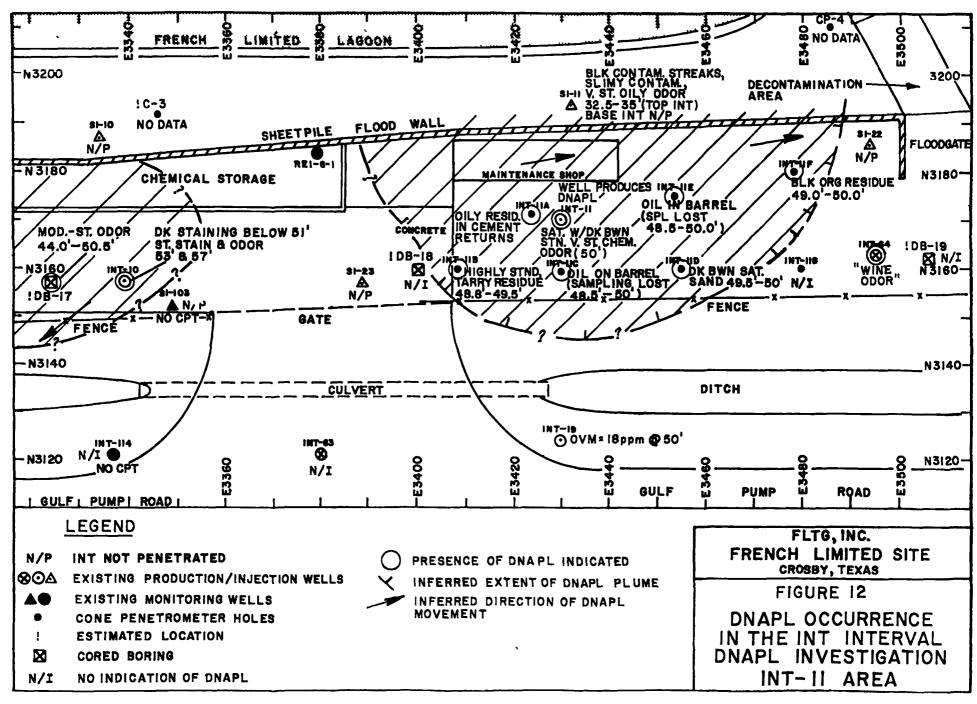


TABLE 1
SUMMARY OF SOIL SAMPLE ANALYSES
\$1-16 AND INT-11 AREAS

WELL	Depth	TOC(ppm)	TPH(ppm)
INT-10	29-30'	BDL	NA
	30-31.5'	BDL	NA
	41-42'	BDL	NA
	50-51'	BDL	NA
	52-53'	BDL	NA
S1-17	11 5-13.5'	BDL	NA
	19-21'	8100	NA
	26-28' (27')	6600	NA
INT-64	15-17'	NA	BDL
	24-26'	NA	BDL
	27-29'	NA	BDL
	31-32'	NA	BDL
	41-43'	NA	BDL
	50-52'	NA	BDL

# TABLE 2

# SAMPLE DESCRIPTIONS S1-16 AND INT-11 AREAS

S1-16A	24' 28'	Water sample. No recovered DNAPL. Water sample. No recovered DNAPL.
\$1-16C	31-32.5' 30.5-32'	Barrel did not open. Red clay at base of sampler.  Moved 2 feet. Lost sample. Mostly sand. Oil remnant in core barrel.
S1-16G	33-33.2'	Sand, gray to blk., m. to c. gr Chemical smell, residual DNAPL. No evidence of mobile product.
	33.2-33.5' 33.5-34.5	Clay, white, silty, with nodules. Chemical smell but no visible staining. Clay, red, mottled with gray silt and f. gr. sand. No DNAPL evident
S1-16H	33-33.5' 33.5-33.8' 33.8-34.5'	Clay, red. No DNAPL. Sand, red, f. gr,. No DNAPL. Clay, red. No DNAPL.
INT-11A	46' 49'	Water sample. No recovered DNAPL Water sample. No recovered DNAPL.
INT-11B	33-34' 34-34.5'	Sand, gray, f. gr., clean Sand, silty, gray, with thin tan-red clay streaks. Black carbonaceous streaks Smelled of chemicals but no reading on OVM.
	48-48.8' 48.8-49.5'	Sand, gray, f. gr., chemical odor but no visible staining.  Sand, highly stained, organics in fine stratifications. Organic is residual, tarry texture, not fluid although can be mobilized when wet.
INT-11C	33.5-35'	Sand, lt gray, f. gr, chemical odor. Spotty, "concretion-like" nodules of organic residual. Noted brownish discoloration in isolated stratified zones which cleave easily. No noticeable grain size difference.
	48.5-50'	Lost soil sample. Oil residue on barrel.
INT-11D	48.5-49.5'	Sand, gray, f. gr., compacted. Chemical odor and light brown chemical staining evident.
	49.5-50'	Sand, dc. bwn., f. gr., organic saturated, above clayey silt to f. gr. sand. Organic material tends to accentuate stratification in f gr sands. Organic staining peters out in underlying silty/clayey zone.
INT-11E	48 5-50'	Lost soil sample. Core barrel had an oily residue.
INT-11F	49-50'	Sand, f. to m. gr. Impregnated with black organic residue material High clay content Sample dense but malleable. No evidence of free liquid organics. Organic appears to be immobile.
	50-50.5'	Clay, dk. bwn., interbedded with silt and f. gr. sand. Organic residual does not penetrate into clay, but some of the silt/f gr. sand streaks are stained.
INT-11G	49-50.5' 51.5-53'	Lost soil sample. No evidence of oil in soil barrel.  Lost soil sample. No evidence of oil in soil barrel

TABLE 3

VOLATILE ORGANIC ANALYSIS - WELL INT-64

INT-11 AREA

Well INT-64	15-17'	24-26'	27-29'	31-32'	41-43'	50-52'
Parameter	ppb	ppb	ppb	ppb	ppb	ppb
Chloromethane	BDL	BDL	BDL	BDL	BDL	BDL
Bromomethane	BDL	BDL	BDL	BDL	BDL	BDL
Vinyl Chloride	BDL	BDL	BDL	BDL	BDL	BDL
Chloroethane	BDL	BDL	BDL	BDL	BDL	BDL
Methylene Chloride	61	130	120	110	*4900	*5100
Acetone	24	12	*1900	<b>*</b> 560	2400	3400
Carbon Disulfide	13	34	BDL	BDL	31000	2800
1,1-Dichloroethene	BDL	BDL	BDL	BDL	BDL	BDL
1,1-Dichloroethane	BDL	BDL	BDL	BDL	BDL	BDL
Trans 1,2-Dichloroethene	BDL	10	43	BDL	BDL	BDL
Chloroform	8	30	220	BDL	830	1000
1,2-Dichloroethane	14	140	1500	260	2800	4400
2-Butanone	BDL	BDL	390	100	BDL	BDL
1,1,1-Trichloroethane	BDL	BDL	BDL	BDL	BDL	BDL
Carbon Tetrachloride	BDL	BDL	BDL	BDL	BDL	BDL
Vınyl Acetate	BDL	BDL	320	75	BDL	BDL
Bromodichloromethane	BDL	BDL	BDL	BDL	BDL	BDL
1,2-Dichloropropane	BDL	BDL	BDL	BDL	BDL	BDL
Cis 1,3-Dichloropropene	BDL	BDL	BDL	BDL	BDL	BDL
Trichloroethene	BDL	BDL	BDL	BDL	BDL	BDL
Dibromochloromethane	BDL	BDL	BDL	BDL	BDL	BDL
1,1,2-Trichloroethane	BDL	BDL	BDL	BDL	BDL	BDL
Benzene	BDL	BDL	BDL	BDL	BDL	BDL
Trans 1,3-Dichloropropene	BDL	BDL	BDL	BDL	BDL	BDL
2-Chloroethylvinylether	BDL	BDL	BDL	BDL	BDL	BDL
Bromoform	BDL	BDL	BDL	BDL	BDL	BDL
4-Methyl 2-Pentanone	BDL	BDL	190	BDL	BDL	BDL
2-Hexanone	BDL	BDL	BDL	BDL	BDL	BDL
Tetrachloroethene	BDL	BDL	BDL	BDL	BDL	BDL
1,1,2,2-Tetrachloroethane	BDL	BDL	BDL	BDL	BDL	BDL
Toluene	BDL	BDL	BDL	BDL	BDL	BDL
Chlorobenzene	BDL	BDL	BDL	BDL	BDL	BDL
Ethylbenzene	BDL	BDL	86	BDL	BDL	BDL
Styrene	BDL	BDL	BDL	BDL	BDL	BDL
Total Xylenes	BDL	BDL	BDL	BDL	BDL	BDL

<sup>\*</sup> Compound Detected in the QC Blank

TABLE 3

VOLATILE ORGANIC ANALYSES - WELL INT-64

INT-11 AREA

Well INT-64	15-17'	24-26'	27-29'	31-32'	41-43'	50-52'	DNAPL (1)
PARAMETER	ppb	ppb	ppb	ppb	ppb	ppb	mg/l
TOC	NA	NA	NA	NA	NA	NA	506
TOX	NA	NA	NA	NA	NA	NA	595
Volatile Organics:							
Chloromethane	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Bromomethane	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Vinyl Chloride	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Chloroethane	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Methylene Chloride	61	130	120	110	4900 (2)	5100 (2)	20,000 (3)
Acetone	24	12	1900 (2)	560 (2)	2400	3400	BDL
Carbon Disulfide	13	34	BDL	BDL	31000	2800	BDL
1,1-Dichloroethene	BDL	BDL	BDL	BDL	BDL	BDL	BDL
1,1-Dichloroethane	BDL	BDL	BDL	BDL	BDL	BDL	20,000 (3)
Trans 1,2-Dichloroethene	BDL	10	43	BDL	BDL	BDL	2,100 (3)
Chloroform	8	30	<b>22</b> 0	BDL	830	1000	810,000
1,2-Dichloroethane	14	140	1500	260	2800	4400	650,000
2-Butanone	BDL	BDL	390	100	BDL	BDL	23,000 (3)
1,1,1-Trichloroethane	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Carbon Tetrachloride	BDL	BDL	BDL	BDL	BDL	BDL	31,000 (3)
Vınyl Acetate	BDL	BDL	320	75	BDL	BDL	150,000
Bromodichloromethane	BDL	BDL	BDL	BDL	BDL	BDL	BDL
1,2-Dichloropropane	BDL	BDL	BDL	BDL	BDL	BDL	BDL
C1s 1,3-D1chloropropene	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Trichloroethene	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Dibromochloromethane	BDL	BDL	BDL	BDL	BDL	BDL	BDL
1,1,2-Trichloroethane	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Benzene	BDL	BDL	BDL	BDL	BDL	BDL	4,400 (3)
Trans 1,3-Dichloropropene	BDL	BDL	BDL	BDL	BDL	BDL	BDL
2-Chloroethylvmylether	BDL	BDL	BDL	BDL	BDL	BDL	NA
Bromoform	BDL	BDL	BDL	BDL	BDL	BDL	BDL
4-Methyl 2-Pentanone	BDL	BDL	190	BDL	BDL	BDL	27,000 (3)
2-Hexanone	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Tetrachloroethene	BDL	BDL	BDL	BDL	BDL	BDL	18,000 (3)
1,1,2,2-Tetrachloroethane	BDL	BDL	BDL	BDL	BDL	BDL	19,000 (3)
Toluene	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Chlorobenzene	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Ethylbenzene	BDL	BDL	<b>8</b> 6	BDL	BDL	BDL	BDL
Styrene	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Total Xylenes	BDL	BDL	BDL	BDL	BDL	BDL	BDL

#### **Notes**

- (1) Sample SPEC-015 collected 4/8/92
- (2) Compound detected in the QC Blank
- (3) Detected at a concentration below the detection limit

Z.ELITRI

TABLE 4

ANALYSIS OF DNAPL SAMPLES FROM INT-11B AND INT-11C

Analyte:	INT-11B	INT-11B(dup)	INT-11C
	mg/kg	mg/kg	mg/kg
Total organic carbon	38,900	39,200	31,700
Total organic halogens	80,300	79,000	112,000
Total petroleum hydrocarbons	97,100		69,500

VOCs	INT-11B (1:1.6)	INT-11B (1.310)	INT-11C (1.1.6)	INT-11C (1.310)
	mg/kg	mg/kg	mg/kg	mg/kg
Vinyl chloride	8J			
Chlorethane	140		5J	
Methylene chloride	210		49	
Acetone				2100J
Carbon disulfide	16		3J	
1,1-Dichloroethene	300		50	
1,1-Dichloroethane	970	520J	380	
1,2-Dichloroethene (total)	7,200	5,000	2,900	
Chloroform	12,000	27,000	8,700	9,200
1,2-Dichloroethane	13,000	12,000	9,100	9,200
2-Butanone			18	,
1,1,1-Trichloroethane	1		4J	;
Carbon tetrachloride	7,000	26,000	5,600	7,500
Trichloroethene	700	700J	480	
1,1 2-Trichloroethane	12		8J	·
Benzene	27		21	
4-Methyl-2-pentanone	54		40	
2-Hexanone	73		43	
Tetrachloroethene	3,700	13,000	3,500	4,500
1,1,2,2-Tetrachloroethane	130		100	'
Toluene	90		51	
Ethylbenzene	87		54	
Xylenes (total)	250		170	
Tentatively Identified:		1	}	
Ethane, pentachloro-	690		340	

SVOC <sub>3</sub>	INT-11B	INT-11C	
	mg/kg	mg/kg	
Hexachloroethane	5,800J	2,200J	
Naphthalene	1,600J	660J	
Hexachlorobutadiene	38,000	17,000	
2-Methylnaphthalene	350J		
Acenaphthene	610J		
Dibenzofuran	360Ј		
Fluorene	(TOO)		
Phenanthrene	1,000J	410J	
Tentatively Identified.			
1,3-Butadiene, pentachloro-	5,900	1	
Aliphatic hydrocarbon	7,300		

Only compounds detected are shown

J = present, but below the detection limit

#### Summary

Total VOCs (inc J)	INT-11B (1 1 6)	INT-11B (1 310)	INT-11C (1 1 6)	INT-11C (1 310)
	mg/kg	mg/kg	mg/kg	mg/kg
	46,591	84,220		32,500

Total SVOCs (inc J)	INT-11B	INT-11C	
	mg/kg	 mg/kg	
	61,520	20,270	

# **BOOKMARK**

Appendix 1

Analytical Reports S1-16 DNAPL sample

Report Prepared for

FLTG , INC.

1024 GULF PUMP ROAD. CROSBY TX 77532

Attention : Ted Davis

TO: 430 VIA: Course

FEB 14 1992

by

**Keystone Lab - Houston** A Division of CHESTER LabNet 8300 Westpark Drive Houston, Texas 77063 (713) 266-6800

Dan Pastalaniec

Project Manager

WORK ORDER

IAH 3/16/92

: H92-01.331

DATE RECEIVED : 30-JAN-1992

PROJECT ID

P.O. NUMBER : SPEC001-01

#### SAMPLE & ANALYSIS SUMMARY

Keystone Sample ID	Client's Sample Number	Date/Time Collected	Sample Matrix
H92-01.331-001 H92-01.331-002		01/30/92 08:45	WATER WATER
Volatiles			
Date analyzed	Dilution   Analyst		
02/07/92	500K   PAA		
Semivolatiles			
Date analyzed	Dilution   Analyst		
02/06/92	100   JDD		
Int Vol BN	Fin Vol EXT date  Ana	alyst  Int Vol A	Fin Vol EXT date
1g	10 ml  01/25/92  BF	PC	1
Surrogate #	Date Prep.		
6226A	01/25/92	<del></del>	

COMMENTS: This sample was done as a waste dilution. There is about 5% free water in the sample and all of the rest appears to be soluble in methylene chloride.

# ORGANICS ANALYSIS DATA SHEET

LABORATORY NAME: KEYSTONE ENV.

3 SAMPLE ID NO.: 910133102

.4PLE MATRIX: WATER

DATA RELEASE AUTHORIZED BY: . !

CASE NO. : ----

INITIAL CALIBRATION DATE: 02/05/92

CONTRACT NO. : --

DATE SAMPLE RECEIVED: 01/30/92

CONCENTRATION: LOW DATAFILE: 4U01331V02B

DATE ANALYZED: 02/07/92 DILUTION FACTOR: 500.00 K

**PECATILES** 

	COMPOUND	DETECTION LIMIT (MILLIGRAM	AMOUNT FOUND MS / LITER )
 CO10	CHLOROMETHANE	5000 U	
	BROMOMETHANE	5000 U	
C020	e : : : : : : : : : : : : : : : : : : :	5000 U	
C025		5000 U	
C030	CHLOROETHANE METHYLENE CHLORIDE ACETONE	2500 U	
CQ35	ACETONE	5000 U	
C040		2500 U	
CO45	1.1-DICHLOROETHENE	2500 U	
C050	1,1-DICHLOROETHANE	2500	2000 J
C053	1,2-DICHLORGETHENE (TOTAL)		2300 J
C040	CHLOROFORM	2500	78000
C065	1,2-DICHLOROETHANE	2500	65000
C110	2-BUTANONE	5000 U	
		2500 U	
	CARBON TETRACHLORIDE	2500	20000
C125	VINYL ACETATE	5000 U	
C130	BROMODICHLOROMETHANE	2500 U	
C140	1,2-DICHLOROPROPANE	2500 U	
C143	CIS-1.3-DICHLOROPROPENE	2500 U	
	TRICHLOROETHENE	2500 U	Loop
	DIBROMOCHLOROMETHANE	2500 U 2500 U	
	1, 1, 2-TRICHLORDETHANE	2500 U	
	BENZENE	2500 U	
C172	TRANS-1, 3-DICHLOROPROPENE	2500 U	
C180	BROMOFORM	2500 U	
C205	4-METHYL-2-PENTANONE	5000 U	
C210	2-HEXANONE	5000 U	
C220	TETRACHLOROETHENE	2500 U	
	1, 1, 2, 2-TETRACHLOROETHANE	2500 U	_
C230	TOLUENE	2500 U	780J
C235	CHLOROBENZENE	2500 U	·
C240	ETHYLBENZENE	2500 V	
<del>-</del> -	STYRENE	2500 U	
	XYLENES (TOTAL)		1900 J

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

072454

## VOA QUANTITATION REPORT

SAMPLE ID.

SPEC0001-01 SI-16

F NAME

4U01331V02B

INST

4000

CLIENT

FRENCH

ANALYST

PVG

DATE INJECTED 02/07/92 17:39:00

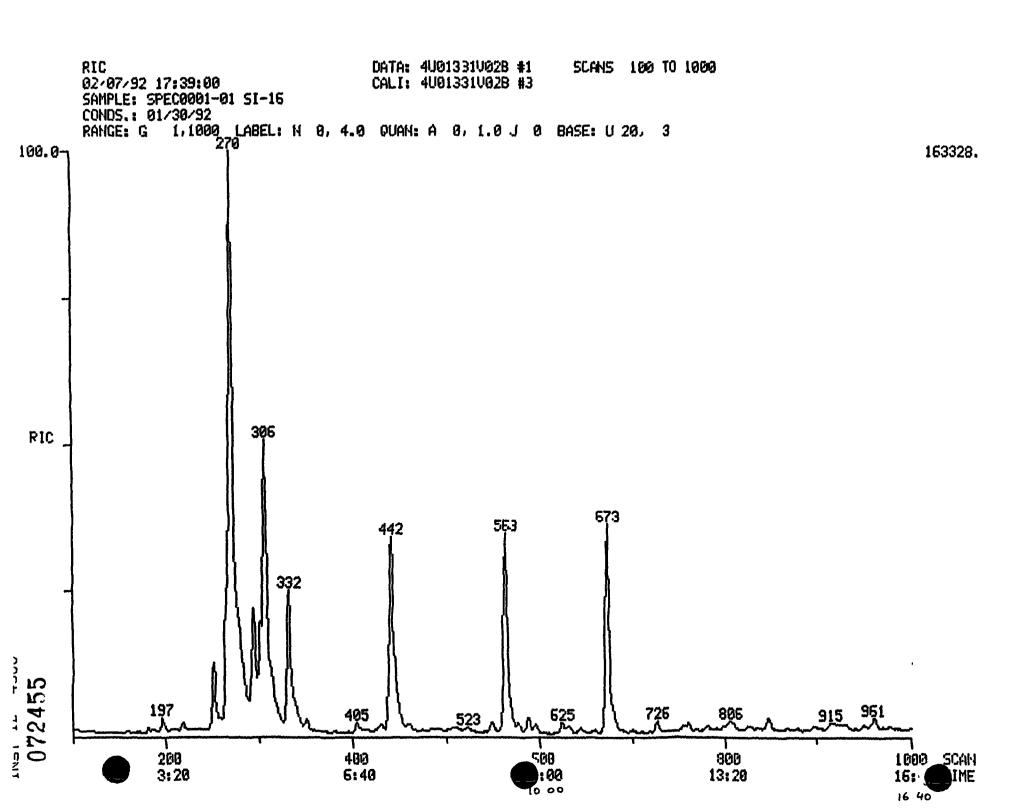
VERIFIED BY

CALIB STD DATE 02/05/92

CORR. FACTOR

500. QQ

				unitima/L		
SCAN#	VOA COMPOUNDS		M/E	AMOUNT	RRT	AREA
265 CI01	BROMOCHLOROMETHANE	 **IS1**	128	25000 06/L	1.000	20114
332 CI10	1,4-DIFLUOROBENZENE	**IS2**	114	25000 QQ/L	1.000	90281
563 CI20	D5-CHLOROBENZENE	**E21#*	117	25000 UQ/L	1.000	87165
.302 CS15	D4-1, 2-DICHLORGETHANE	**5U1**	65	127 )%	1. 140	50050
442 C505	D8-TOLUENE	**SU2**	<b>78</b>	106 %	1. 331	104401
673 C510	P-BROMOFLUOROBENZENE	**EU3**	95	104 %	1. 175	60747
219 C050	1,1-DICHLOROETHANE		63	1955 NG/L	0. 826	71 <del>5</del> 8
197 C053	1, 2-DICHLOROETHENE (TOTA	AL)	96	2335 BQ/L	0. 743	3123
270 CQ60	CHLOROFORM		83	78482   UÇ/L	1.019	296925
307 CO65	1,2-DICHLORGETHANE		62	65300 UC/L	1. 158	152113
294 C120	CARBON TETRACHLORIDE		117	19822 (UB/L	0. 886	46221
588 C250	XYLENES (TOTAL)		106	1855\0C/L	1.044	<b>6550</b>



SAMPLE NUMBER VBLK

072456

#### ORGANICS ANALYSIS DATA SHEET

TRATORY NAME: KEYSTONE ENV.

SAMPLE ID NO.: 910133101

SAMPLE MATRIX: WATER

DATA RELEASE AUTHORIZED BY:.

CASE NO. : ----

INITIAL CALIBRATION DATE: 02/05/92

CONTRACT NO. : --

DATE SAMPLE RECEIVED: ----

DEATILES

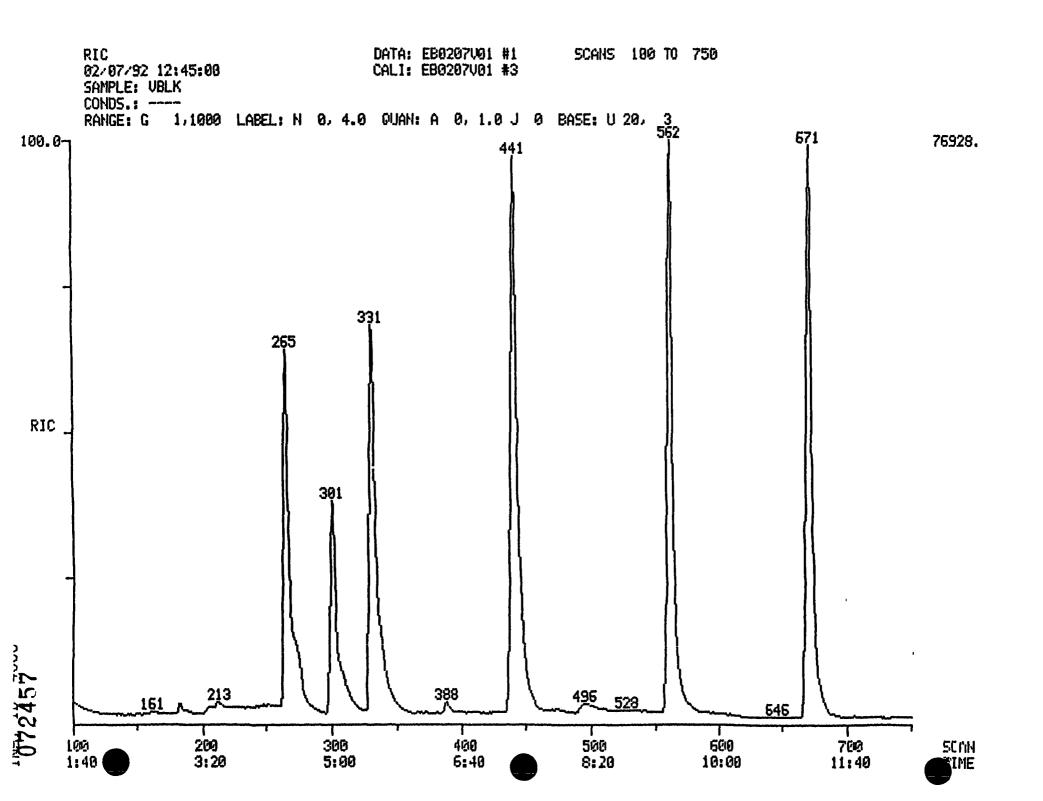
CONCENTRATION: LOW DATAFILE: EB0207V01

DATE ANALYZED: 02/07/92 DILUTION FACTOR: 1.00

CO10 CHLOROMETHANE CO15 BROMOMETHANE CO20 VINYL CHLORIDE CO20 VINYL CHLORIDE CO20 CHLOROETHANE CO30 METHYLENE CHLORIDE CO30 METHYLENE CHLORIDE CO40 CARBON DISULFIDE CO40 CARBON DISULFIDE CO45 1, 1-DICHLOROETHENE CO50 1, 1-DICHLOROETHENE CO50 1, 1-DICHLOROETHENE CO50 1, 2-DICHLOROETHENE CO50 1, 2-DICHLOROETHANE CO50 1, 2-DICHLOROETHANE CO50 1, 2-DICHLOROETHANE CO50 1, 2-DICHLOROETHANE CO10 CARBON TETRACHLORIDE CO10 CARBON TETRACHLORIDE CO10 CARBON TETRACHLORIDE CO10 CARBON TETRACHLORIDE CO10 BROMODICHLOROMETHANE CO10 TICHLOROETHANE CO10 TRICHLOROETHANE CO10 TRICHLOROETHA		COMPOUND	DETECTION AMOUNT LIMIT FOUND (MICROGRAMS / LITER )			
COAS 1.2-DICHLOROETHANE 10 U C110 2-BUTANONE 10 U C115 1.1.1-TRICHLOROETHANE 5 U C120 CARBON TETRACHLORIDE 5 U C125 VINYL ACETATE 10 U C130 BROMODICHLOROMETHANE 5 U C140 1.2-DICHLOROPROPANE 5 U C141 CIS-1.3-DICHLOROPROPENE 5 U C150 TRICHLOROETHANE 5 U C150 TRICHLOROETHANE 5 U C151 DIBROMOCHLOROMETHANE 5 U C152 DIBROMOCHLOROMETHANE 5 U C164 1.1.2-TRICHLOROETHANE 5 U C165 BENZENE 5 U C172 TRANS-1.3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C	010 CHLOROMETHANE	10 U			
COAS 1.2-DICHLOROETHANE 10 U C110 2-BUTANONE 10 U C115 1.1.1-TRICHLOROETHANE 5 U C120 CARBON TETRACHLORIDE 5 U C125 VINYL ACETATE 10 U C130 BROMODICHLOROMETHANE 5 U C140 1.2-DICHLOROPROPANE 5 U C141 CIS-1.3-DICHLOROPROPENE 5 U C150 TRICHLOROETHANE 5 U C150 TRICHLOROETHANE 5 U C151 DIBROMOCHLOROMETHANE 5 U C152 DIBROMOCHLOROMETHANE 5 U C164 1.1.2-TRICHLOROETHANE 5 U C165 BENZENE 5 U C172 TRANS-1.3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C	)15 BROMOMETHANE	10 U			
COAS 1.2-DICHLOROETHANE 10 U C110 2-BUTANONE 10 U C115 1.1.1-TRICHLOROETHANE 5 U C120 CARBON TETRACHLORIDE 5 U C125 VINYL ACETATE 10 U C130 BROMODICHLOROMETHANE 5 U C140 1.2-DICHLOROPROPANE 5 U C141 CIS-1.3-DICHLOROPROPENE 5 U C150 TRICHLOROETHANE 5 U C150 TRICHLOROETHANE 5 U C151 DIBROMOCHLOROMETHANE 5 U C152 DIBROMOCHLOROMETHANE 5 U C164 1.1.2-TRICHLOROETHANE 5 U C165 BENZENE 5 U C172 TRANS-1.3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C	20 VINYL CHLORIDE	10 U			
COAS 1.2-DICHLOROETHANE 10 U C110 2-BUTANONE 10 U C115 1.1.1-TRICHLOROETHANE 5 U C120 CARBON TETRACHLORIDE 5 U C125 VINYL ACETATE 10 U C130 BROMODICHLOROMETHANE 5 U C140 1.2-DICHLOROPROPANE 5 U C141 CIS-1.3-DICHLOROPROPENE 5 U C150 TRICHLOROETHANE 5 U C150 TRICHLOROETHANE 5 U C151 DIBROMOCHLOROMETHANE 5 U C152 DIBROMOCHLOROMETHANE 5 U C164 1.1.2-TRICHLOROETHANE 5 U C165 BENZENE 5 U C172 TRANS-1.3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C	25 CHLOROETHANE	10 U			
COAS 1.2-DICHLOROETHANE 10 U C110 2-BUTANONE 10 U C115 1.1.1-TRICHLOROETHANE 5 U C120 CARBON TETRACHLORIDE 5 U C125 VINYL ACETATE 10 U C130 BROMODICHLOROMETHANE 5 U C140 1.2-DICHLOROPROPANE 5 U C141 CIS-1.3-DICHLOROPROPENE 5 U C150 TRICHLOROETHANE 5 U C150 TRICHLOROETHANE 5 U C151 DIBROMOCHLOROMETHANE 5 U C152 DIBROMOCHLOROMETHANE 5 U C164 1.1.2-TRICHLOROETHANE 5 U C165 BENZENE 5 U C172 TRANS-1.3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C	330 METHYLENE CHLORIDE	5 U			
COAS 1.2-DICHLOROETHANE 10 U C110 2-BUTANONE 10 U C115 1.1.1-TRICHLOROETHANE 5 U C120 CARBON TETRACHLORIDE 5 U C125 VINYL ACETATE 10 U C130 BROMODICHLOROMETHANE 5 U C140 1.2-DICHLOROPROPANE 5 U C141 CIS-1.3-DICHLOROPROPENE 5 U C150 TRICHLOROETHANE 5 U C150 TRICHLOROETHANE 5 U C151 DIBROMOCHLOROMETHANE 5 U C152 DIBROMOCHLOROMETHANE 5 U C164 1.1.2-TRICHLOROETHANE 5 U C165 BENZENE 5 U C172 TRANS-1.3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C	035 ACETONE	10	4 J		
COAS 1.2-DICHLOROETHANE 10 U C110 2-BUTANONE 10 U C115 1.1.1-TRICHLOROETHANE 5 U C120 CARBON TETRACHLORIDE 5 U C125 VINYL ACETATE 10 U C130 BROMODICHLOROMETHANE 5 U C140 1.2-DICHLOROPROPANE 5 U C141 CIS-1.3-DICHLOROPROPENE 5 U C150 TRICHLOROETHANE 5 U C150 TRICHLOROETHANE 5 U C151 DIBROMOCHLOROMETHANE 5 U C152 DIBROMOCHLOROMETHANE 5 U C164 1.1.2-TRICHLOROETHANE 5 U C165 BENZENE 5 U C172 TRANS-1.3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C	040 CARBON DISULFIDE	5 ປ			
COAS 1.2-DICHLOROETHANE 10 U C110 2-BUTANONE 10 U C115 1.1.1-TRICHLOROETHANE 5 U C120 CARBON TETRACHLORIDE 5 U C125 VINYL ACETATE 10 U C130 BROMODICHLOROMETHANE 5 U C140 1.2-DICHLOROPROPANE 5 U C141 CIS-1.3-DICHLOROPROPENE 5 U C150 TRICHLOROETHANE 5 U C150 TRICHLOROETHANE 5 U C151 DIBROMOCHLOROMETHANE 5 U C152 DIBROMOCHLOROMETHANE 5 U C164 1.1.2-TRICHLOROETHANE 5 U C165 BENZENE 5 U C172 TRANS-1.3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C	045 1,1-DICHLOROETHENE	5 U			
COAS 1.2-DICHLOROETHANE 10 U C110 2-BUTANONE 10 U C115 1.1.1-TRICHLOROETHANE 5 U C120 CARBON TETRACHLORIDE 5 U C125 VINYL ACETATE 10 U C130 BROMODICHLOROMETHANE 5 U C140 1.2-DICHLOROPROPANE 5 U C141 CIS-1.3-DICHLOROPROPENE 5 U C150 TRICHLOROETHANE 5 U C150 TRICHLOROETHANE 5 U C151 DIBROMOCHLOROMETHANE 5 U C152 DIBROMOCHLOROMETHANE 5 U C164 1.1.2-TRICHLOROETHANE 5 U C165 BENZENE 5 U C172 TRANS-1.3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C	050 1,1-DICHLOROETHANE	5 U			
COAS 1.2-DICHLOROETHANE 10 U C110 2-BUTANONE 10 U C115 1.1.1-TRICHLOROETHANE 5 U C120 CARBON TETRACHLORIDE 5 U C125 VINYL ACETATE 10 U C130 BROMODICHLOROMETHANE 5 U C140 1.2-DICHLOROPROPANE 5 U C141 CIS-1.3-DICHLOROPROPENE 5 U C150 TRICHLOROETHANE 5 U C150 TRICHLOROETHANE 5 U C151 DIBROMOCHLOROMETHANE 5 U C152 DIBROMOCHLOROMETHANE 5 U C164 1.1.2-TRICHLOROETHANE 5 U C165 BENZENE 5 U C172 TRANS-1.3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C	053 1.2-DICHLOROETHENE (TOTAL)	5 U			
COAS 1.2-DICHLOROETHANE 10 U C110 2-BUTANONE 10 U C115 1.1.1-TRICHLOROETHANE 5 U C120 CARBON TETRACHLORIDE 5 U C125 VINYL ACETATE 10 U C130 BROMODICHLOROMETHANE 5 U C140 1.2-DICHLOROPROPANE 5 U C141 CIS-1.3-DICHLOROPROPENE 5 U C150 TRICHLOROETHANE 5 U C150 TRICHLOROETHANE 5 U C151 DIBROMOCHLOROMETHANE 5 U C152 DIBROMOCHLOROMETHANE 5 U C164 1.1.2-TRICHLOROETHANE 5 U C165 BENZENE 5 U C172 TRANS-1.3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	_ CC	060 CHLOROFORM	5 U			
C115 1.1.1—TRICHLOROETHANE 5 U C120 CARBON TETRACHLORIDE 5 U C125 VINYL ACETATE 10 U C130 BROMODICHLOROMETHANE 5 U C140 1.2—DICHLOROPROPANE 5 U C143 CIS—1.3—DICHLOROPROPENE 5 U C150 TRICHLOROETHENE 5 U C150 DIBROMOCHLOROMETHANE 5 U C155 DIBROMOCHLOROMETHANE 5 U C160 1.1.2—TRICHLOROETHANE 5 U C165 BENZENE 5 U C172 TRANS—1.3—DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4—METHYL—2—PENTANONE 10 U C210 2—HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1.1.2.2—TETRACHLOROETHANE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	CC	065 1.2-DICHLOROETHANE	5 U			
C115 1.1.1—TRICHLOROETHANE 5 U C120 CARBON TETRACHLORIDE 5 U C125 VINYL ACETATE 10 U C130 BROMODICHLOROMETHANE 5 U C140 1.2—DICHLOROPROPANE 5 U C143 CIS—1.3—DICHLOROPROPENE 5 U C150 TRICHLOROETHENE 5 U C150 DIBROMOCHLOROMETHANE 5 U C155 DIBROMOCHLOROMETHANE 5 U C160 1.1.2—TRICHLOROETHANE 5 U C165 BENZENE 5 U C172 TRANS—1.3—DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4—METHYL—2—PENTANONE 10 U C210 2—HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1.1.2.2—TETRACHLOROETHANE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C	10 2-BUTANONE	10 U			
C130 BROMODICHLOROMETHANE 5 U C140 1,2-DICHLOROPROPANE 5 U C143 CIS-1,3-DICHLOROPROPENE 5 U C150 TRICHLOROETHENE 5 U C155 DIBROMOCHLOROMETHANE 5 U C160 1,1,2-TRICHLOROETHANE 5 U C161 BENZENE 5 U C172 TRANS-1,3-DICHLOROPROPENE 5 U C172 TRANS-1,3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1,1,2,2-TETRACHLOROETHANE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C	15 1,1,1-TRICHLOROETHANE	5 U			
C130 BROMODICHLOROMETHANE 5 U C140 1,2-DICHLOROPROPANE 5 U C143 CIS-1,3-DICHLOROPROPENE 5 U C150 TRICHLOROETHENE 5 U C155 DIBROMOCHLOROMETHANE 5 U C160 1,1,2-TRICHLOROETHANE 5 U C161 BENZENE 5 U C172 TRANS-1,3-DICHLOROPROPENE 5 U C172 TRANS-1,3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1,1,2,2-TETRACHLOROETHANE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C	20 CARBON TETRACHLORIDE	5 U			
C130 BROMODICHLOROMETHANE 5 U C140 1,2-DICHLOROPROPANE 5 U C143 CIS-1,3-DICHLOROPROPENE 5 U C150 TRICHLOROETHENE 5 U C155 DIBROMOCHLOROMETHANE 5 U C160 1,1,2-TRICHLOROETHANE 5 U C161 BENZENE 5 U C172 TRANS-1,3-DICHLOROPROPENE 5 U C172 TRANS-1,3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1,1,2,2-TETRACHLOROETHANE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	Ci	25 VINYL ACETATE	10 U			
C172 TRANS-1, 3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1, 1, 2, 2-TETRACHLOROETHANE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C1	30 BROMODICHLOROMETHANE	5 U			
C172 TRANS-1, 3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1, 1, 2, 2-TETRACHLOROETHANE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C1	40 1,2-DICHLOROPROPANE	5 U			
C172 TRANS-1, 3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1, 1, 2, 2-TETRACHLOROETHANE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	CI	43 CIS-1.3-DICHLOROPROPENE	5 V			
C172 TRANS-1, 3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1, 1, 2, 2-TETRACHLOROETHANE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	CI	50 TRICHLORDETHENE	5 U			
C172 TRANS-1, 3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1, 1, 2, 2-TETRACHLOROETHANE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	CI	55 DIBROMOCHLOROMETHANE	5 U			
C172 TRANS-1, 3-DICHLOROPROPENE 5 U C180 BROMOFORM 5 U C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1, 1, 2, 2-TETRACHLOROETHANE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C1	.60 1,1,2-TRICHLORDETHANE	5 U			
C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1,1,2,2-TETRACHLOROETHANE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C1	.65 BENZENE	5 U			
C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1,1,2,2-TETRACHLOROETHANE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	CI	72 TRANS-1,3-DICHLOROPROPENE	5 U			
C205 4-METHYL-2-PENTANONE 10 U C210 2-HEXANONE 10 U C220 TETRACHLOROETHENE 5 U C225 1,1,2,2-TETRACHLOROETHANE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	C1	80 BROMOFORM	5 U			
C210 2-HEXANDNE 10 U C220 TETRACHLOROETHENE 5 U C225 1,1,2,2-TETRACHLOROETHANE 5 U C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	Ca	IOS A-METUVI -9-BENTANIONE	10 11			
C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U	CZ	10 2-HEXANONE	10 U			
C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U		20 TETRACHLOROETHENE	5 U			
C230 TOLUENE 5 U C235 CHLOROBENZENE 5 U C240 ETHYLBENZENE 5 U C245 STYRENE 5 U		25 1, 1, 2, 2-TETRACHLOROETHANE	5 U			
C243 STRENE S V		30 TOLUENE	ט כ			
C243 STRENE S V		35 CHLOROBENZENE				
C243 STRENE S V		40 ETHYLBENZENE	<del>-</del>			
C250 XYLENES (TDTAL) 5 U		AN SITE CH				
	CE	50 XYLENES (TOTAL)	5 ບ			

U = UNDETECTED AT THE LISTED DETECTION LIMIT

<sup>-</sup> COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT



072458

## VOA QUANTITATION REPORT

SAMPLE ID.

VBLK

ENAME

EB0207V01

INST

4000

CLIENT

FRENCH

ANALYST

PVG

DATE INJECTED

02/07/92 12:45:00

VERIFIED BY

CALIB STD DATE 02/05/92

CORR. FACTOR

1.00

SCAN#	VOA COMPOUNDS		M/E	AMOUNT	RRT	AREA
265 CI01	BROMOCHLOROMETHANE	**IS1**	128	50 UG/L	1.000	24575
331 CI10	1,4-DIFLUOROBENZENE	**IS2**	114	50 UG/L	1.000	119864
542 CI20	D5-CHLOROBENZENE	**IS3**	117	50 UG/L	1.000	117539
301 CS15	D4-1, 2-DICHLOROETHANE	**SU1**	<b>65</b>	113 %	1. 136	54624
441 CS05	D8-TOLUENE	**SU2**	<del>98</del>	102 %	1. 332	133867
671 CS10	P-BROMOFLUOROBENZENE	**SU3**	<b>9</b> 5	100 %	1. 194	78993
161 CO35	ACETONE		43	4 UG/L	0. 608	1272

#### ORGANICS ANALYSIS DATA SHEET -

DRATORY NAME: KEYSTONE

6 SAMPLE ID NO.: 910133102

SAMPLE MATRIX. SOIL

DATA RELEASE AUTHORIZED BY:

CASE NO. ---

QC REPORT NO. . 01/27/92

CONTRACT NO . ---

DATE SAMPLE RECEIVED: ... 1/30/93

SEMIVOLATILES

CONCENTRATION: LOW DATE EXTRACTED: ..!!!!!

DATE ANALYZED: 02/06/92

DATAFILE: 9U01331C02

C555 2,4-DINITROPHENOL 5000 U C560 4-NITROPHENOL 5000 U C565 DIBENZOFURAN 1000 230 C C570 2,4-DINITROTOLUENE 1000 U			DETECTION LIMIT (MILLIGRAMS)	AMOUNT FOUND / KG)
C330 2-CHLQROPHENOL 1000 U C335 1,3-D1CHLQROBENZENE 1000 U C345 BENZYL ALCOHOL 1000 U C356 1,2-D1CHLQROBENZENE 1000 U C357 2,2'-D1CHLQROBENZENE 1000 U C357 2,2'-GXYBIS(1-CHLQROPROPANE) 1000 U C357 2,2'-GXYBIS(1-CHLQROPROPANE) 1000 U C357 4,4-METHYLPHENOL 1000 U C357 N-NITROSODIPROPYLAMINE 1000 U C370 N-NITROSODIPROPYLAMINE 1000 U C410 NITROBENZENE 1000 U C415 ISOPHORONE 1000 U C420 2,NITROPHENOL 1000 U C425 2,4-DIMETHYLPHENOL 1000 U C430 BENZOIC ACID 5000 U C430 BENZOIC ACID 5000 U C440 2,4-DICHLQROPHENOL 1000 U C445 1,2,4-TRICHLQROBENZENE 1000 U C445 1,2,4-TRICHLQROBENZENE 1000 U C450 NAPHTHALENE 1000 U C465 P-CHLQRO-TADIENE 1000 U C465 P-CHLQRO-TADIENE 1000 U C465 P-CHLQRO-TADIENE 1000 U C510 PEXACHLOROCYCLOPENTADIENE 1000 U C520 2,4,5-TRICHLOROPHENOL 1000 U C520 2,4,5-TRICHLOROPHENOL 5000 U C520 2,4-DINITROPHENOL 5000 U		PHENOL	1000 U	
C330 2-CHLOROPHENOL 1000 U C335 1,3-DICHLOROBENZENE 1000 U C345 BENZYL ALCOHOL 1000 U C350 1,2-DICHLOROBENZENE 1000 U C357 2,2'-DICHLOROBENZENE 1000 U C357 2,2'-DICHLOROBENZENE 1000 U C357 2,2'-DICHLOROBENZENE 1000 U C357 2,2'-DICHLOROPROPANE) 1000 U C357 4-METHYLPHENOL 1000 U C357 N-NITROSODIPROPYLAMINE 1000 U C370 N-NITROSODIPROPYLAMINE 1000 U C410 NITROBENZENE 1000 U C411 ISOPHORONE 1000 U C412 2-NITROPHENOL 1000 U C420 2-NITROPHENOL 1000 U C430 BENZOIC ACID 5000 U C430 BENZOIC ACID 5000 U C440 2,4-DICHLOROPHENOL 1000 U C444 1,2,4-TRICHLOROBENZENE 1000 U C455 4-CHLOROBENZENE 1000 U C455 4-CHLOROBITADIENE 1000 U C465 P-CHLORO-M-CRESOL 1000 U C465 P-CHLORO-M-CRESOL 1000 U C515 2,4,6-TRICHLOROPHENOL 1000 U C520 2,4,5-TRICHLOROPHENOL 1000 U C520 2,1TROANILINE 1000 U C530 DIMETHYL PHTHALATE 1000 U C540 ACENAPHTHYLENE 1000 U C550 2,4-DINITROPHENOL 5000 U C550 DIBENZOFURAN 1000 U C550 DIBENZOFURAN 1000 U C550 DIBENZOFURAN 1000 U	C325	BIS(2-CHLORDETHYL)ETHER	1000 U	
C340 1, 4-DICHLOROBENZENE 1000 U C345 BENZYL ALCOHOL 1000 U C350 1, 2-DICHLOROBENZENE 1000 U C357 2, 2'-OXYBIS(1-CHLOROPROPANE) 1000 U C357 2, 2'-OXYBIS(1-CHLOROPROPANE) 1000 U C358 4-METHYLPHENOL 1000 U C370 N-NITROSODIPROPYLAMINE 1000 U C370 N-NITROSODIPROPYLAMINE 1000 U C371 HEXACHLOROETHANE 1000 U C410 NITROBENZENE 1000 U C411 ISOPHORONE 1000 U C420 2-NITROPHENOL 1000 U C425 2, 4-DIMETHYLPHENOL 1000 U C426 BENZOIC ACID 5000 U C430 BIS(2-CHLOROETHOXY)METHANE 1000 U C445 1, 2, 4-TRICHLOROBENZENE 1000 U C445 1, 2, 4-TRICHLOROBENZENE 1000 U C450 NAPHTHALENE 1000 U C450 HEXACHLOROBUTADIENE 1000 U C450 P-CHLORO-M-CRESOL 1000 U C450 2-METHYLNAPHTHALENE 1000 U C550 2, 4, 5-TRICHLOROPHENOL 1000 U C551 2, 4, 6-TRICHLOROPHENOL 5000 U C552 2-CHLORO-M-CRESOL 1000 U C553 DIMETHYL PHTHALATE 1000 U C554 3-NITROANILINE 1000 U C555 2-CHLORONAPHTHALENE 1000 U C556 3-NITROANILINE 5000 U C556 3-NITROANILINE 5000 U C5570 2, 4-DINITROPHENOL 5000 U	C330	2-CHLOROPHENOL	1000 U	
C340 1, 4-DICHLOROBENZENE 1000 U C345 BENZYL ALCOHOL 1000 U C350 1, 2-DICHLOROBENZENE 1000 U C357 2, 2'-OXYBIS(1-CHLOROPROPANE) 1000 U C357 2, 2'-OXYBIS(1-CHLOROPROPANE) 1000 U C358 4-METHYLPHENOL 1000 U C370 N-NITROSODIPROPYLAMINE 1000 U C370 N-NITROSODIPROPYLAMINE 1000 U C371 HEXACHLOROETHANE 1000 U C410 NITROBENZENE 1000 U C411 ISOPHORONE 1000 U C420 2-NITROPHENOL 1000 U C425 2, 4-DIMETHYLPHENOL 1000 U C426 BENZOIC ACID 5000 U C430 BIS(2-CHLOROETHOXY)METHANE 1000 U C445 1, 2, 4-TRICHLOROBENZENE 1000 U C445 1, 2, 4-TRICHLOROBENZENE 1000 U C450 NAPHTHALENE 1000 U C450 HEXACHLOROBUTADIENE 1000 U C450 P-CHLORO-M-CRESOL 1000 U C450 2-METHYLNAPHTHALENE 1000 U C550 2, 4, 5-TRICHLOROPHENOL 1000 U C551 2, 4, 6-TRICHLOROPHENOL 5000 U C552 2-CHLORO-M-CRESOL 1000 U C553 DIMETHYL PHTHALATE 1000 U C554 3-NITROANILINE 1000 U C555 2-CHLORONAPHTHALENE 1000 U C556 3-NITROANILINE 5000 U C556 3-NITROANILINE 5000 U C5570 2, 4-DINITROPHENOL 5000 U	C335	1,3-DICHLOROBENZENE	1000 U	
C350 1,2-DICHLOROBENZENE 1000 U C357 2,2'-QXYBIS(1-CHLOROPROPANE) 1000 U C365 4-METHYLPHENOL 1000 U C365 4-METHYLPHENOL 1000 U C370 N-NITROSODIPROPYLAMINE 1000 U C375 HEXACHLOROETHANE 1000 U C410 NITROBENZENE 1000 U C411 ISOPHORONE 1000 U C420 2-NITROPHENOL 1000 U C420 2-NITROPHENOL 1000 U C430 BENZOIC ACID 5000 U C435 BIS(2-CHLOROETHOXY)METHANE 1000 U C440 2, 4-DICHLOROPHENOL 1000 U C440 2, 4-DICHLOROBENZENE 1000 U C455 4-CHLOROBILINE 1000 U C455 4-CHLOROBILINE 1000 U C465 P-CHLORO-M-CRESOL 1000 U C465 P-CHLORO-M-CRESOL 1000 U C510 C510 C510 C510 C510 C510 C510 C510	C340	1,4-DICHLOROBENZENE	1000 U	
C355 2-METHYLPHENOL 1000 U C357 2,2'-OXYBIS(1-CHLOROPROPANE) 1000 U C365 4-METHYLPHENOL 1000 U C370 N-NITROSODIPROPYLAMINE 1000 U C375 HEXACHLOROETHANE 1000 U C410 NITROBENZENE 1000 U C415 ISOPHORONE 1000 U C420 2-NITROPHENOL 1000 U C425 2,4-DIMETHYLPHENOL 1000 U C435 BIS(2-CHLOROETHANE 1000 U C435 BIS(2-CHLOROETHOXY)METHANE 1000 U C440 2,4-DICHLOROPHENOL 1000 U C441 1,2,4-TRICHLOROBENZENE 1000 U C450 NAPHTHALENE 1000 U C450 HEXACHLOROBUTADIENE 1000 U C465 P-CHLOROANILINE 1000 U C465 P-CHLOROM-M-CRESOL 1000 U C470 2-METHYLNAPHTHALENE 1000 U C510 HEXACHLOROSYLOPENTADIENE 1000 U C510 Z,4,5-TRICHLOROPHENOL 1000 U C520 Z,4,5-TRICHLOROPHENOL 1000 U C520 Z-A,5-TRICHLOROPHENOL 1000 U C520 Z-CHLORONAPHTHALENE 1000 U C530 Z-NITROANILINE 1000 U C530 Z-NITROANILINE 1000 U C530 Z-NITROANILINE 1000 U C550 ACENAPHTHENE 1000 U	C345		1000 U	
C355 2-METHYLPHENOL 1000 U C357 2,2'-OXYBIS(1-CHLOROPROPANE) 1000 U C365 4-METHYLPHENOL 1000 U C370 N-NITROSODIPROPYLAMINE 1000 U C375 HEXACHLOROETHANE 1000 U C410 NITROBENZENE 1000 U C415 ISOPHORONE 1000 U C420 2-NITROPHENOL 1000 U C425 2,4-DIMETHYLPHENOL 1000 U C435 BIS(2-CHLOROETHANE 1000 U C435 BIS(2-CHLOROETHOXY)METHANE 1000 U C440 2,4-DICHLOROPHENOL 1000 U C441 1,2,4-TRICHLOROBENZENE 1000 U C450 NAPHTHALENE 1000 U C450 HEXACHLOROBUTADIENE 1000 U C465 P-CHLOROANILINE 1000 U C465 P-CHLOROM-M-CRESOL 1000 U C470 2-METHYLNAPHTHALENE 1000 U C510 HEXACHLOROSYLOPENTADIENE 1000 U C510 Z,4,5-TRICHLOROPHENOL 1000 U C520 Z,4,5-TRICHLOROPHENOL 1000 U C520 Z-A,5-TRICHLOROPHENOL 1000 U C520 Z-CHLORONAPHTHALENE 1000 U C530 Z-NITROANILINE 1000 U C530 Z-NITROANILINE 1000 U C530 Z-NITROANILINE 1000 U C550 ACENAPHTHENE 1000 U	C350	1,2-DICHLOROBENZENE	1000 U	
C365 4-METHYLPHENOL 1000 U C370 N-NITROSODIPROPYLAMINE 1000 U C375 HEXACHLOROETHANE 1000 U C410 NITROBENZENE 1000 U C415 ISOPHORONE 1000 U C420 2-NITROPHENOL 1000 U C425 2, 4-DIMETHYLPHENOL 1000 U C430 BENZOIC ACID 5000 U C435 BIS(2-CHLOROETHOXY)METHANE 1000 U C445 1, 2, 4-DICHLOROPHENOL 1000 U C445 1, 2, 4-TICHLOROBENZENE 1000 U C445 1, 2, 4-TICHLOROBENZENE 1000 U C455 4-CHLOROBUTADIENE 1000 U C455 4-CHLOROBUTADIENE 1000 U C465 P-CHLORO-M-CRESOL 1000 U C510 HEXACHLOROBUTADIENE 1000 U C510 HEXACHLOROBUTADIENE 1000 U C515 2, 4, 6-TRICHLOROPHENOL 1000 U C525 2-CHLOROMAPHTHALENE 1000 U C525 2-CHLOROMAP	C355	2-METHYLPHENOL	1000 U	
C370 N-NITROSODIPROPYLAMINE 1000 U C375 HEXACHLOROETHANE 1000 U C410 NITROBENZENE 1000 U C415 ISOPHORONE 1000 U C420 2-NITROPHENDL 1000 U C425 2, 4-DIMETHYLPHENDL 1000 U C436 BENZOIC ACID 5000 U C437 BIS(2-CHLOROETHOXY)METHANE 1000 U C430 BENZOIC ACID 1000 U C431 I, 2, 4-TRICHLOROBENZENE 1000 U C445 1, 2, 4-TRICHLOROBENZENE 1000 U C450 NAPHTHALENE 1000 U C451 4-CHLOROBUTADIENE 1000 U C452 4-CHLOROBUTADIENE 1000 U C453 P-CHLORO-M-CRESOL 1000 U C454 P-CHLORO-M-CRESOL 1000 U C545 2-METHYLNAPHTHALENE 1000 U C510 EXACHLOROBYTADIENE 1000 U C510 Z, 4, 5-TRICHLOROPHENOL 1000 U C520 Z, 4, 5-TRICHLOROPHENOL 5000 U C520 Z-NITROANILINE 5000 U C530 Z-NITROANILINE 5000 U C530 Z-NITROANILINE 5000 U C530 Z-NITROANILINE 5000 U C530 ACENAPHTHYLENE 1000 . 170 C C540 ACENAPHTHENE 1000 . 740 C C550 ACENAPHTHENE 1000 U C555 Z, 4-DINITROPHENOL 5000 U C556 DIBENZOFURAN 1000 . 230 C C570 Z, 4-DINITROTOLUENE 1000 U	C357	2,2'-OXYBIS(1-CHLOROPROPANE)	1000 U	
C375 HEXACHLOROETHANE 1000 U C410 NITROBENZENE 1000 U C415 ISOPHORONE 1000 U C420 2-NITROPHENDL 1000 U C422 2,4-DIMETHYLPHENDL 1000 U C425 2,4-DIMETHYLPHENDL 1000 U C430 BENZOIC ACID 5000 U C430 BIS(2-CHLOROETHOXY)METHANE 1000 U C445 1,2,4-TRICHLOROPHENDL 1000 U C445 1,2,4-TRICHLOROBENZENE 1000 U C450 NAPHTHALENE 1000 U C455 4-CHLOROANILINE 1000 U C455 4-CHLOROBUTADIENE 1000 U C465 P-CHLORO-M-CRESOL 1000 U C470 2-METHYLNAPHTHALENE 1000 U C510 HEXACHLOROCYCLOPENTADIENE 1000 U C510 HEXACHLOROCYCLOPENTADIENE 1000 U C520 2,4,5-TRICHLOROPHENDL 1000 U C525 2-CHLORONAPHTHALENE 1000 U C535 DIMETHYL PHTHALATE 1000 U C535 DIMETHYL PHTHALATE 1000 U C540 ACENAPHTHYLENE 1000  170 C C550 ACENAPHTHENE 1000  740 C C550 ACENAPHTHENE 1000 U C555 2,4-DINITROPHENOL 5000 U C555 2,4-DINITROPHENOL 5000 U C555 DIBENZOFURAN 1000  230 C C570 2,4-DINITROTOLUENE 1000 U			1000 U	
C410 NITROBENZENE 1000 U C415 ISOPHORONE 1000 U C420 2-NITROPHENDL 1000 U C425 2, 4-DIMETHYLPHENDL 1000 U C425 2, 4-DIMETHYLPHENDL 1000 U C435 BIS(2-CHLOROETHOXY)METHANE 1000 U C440 2, 4-DICHLOROPHENDL 1000 U C441 1, 2, 4-TRICHLOROBENZENE 1000 U C445 1, 2, 4-TRICHLOROBENZENE 1000 U C450 NAPHTHALENE 1000 U C455 4-CHLOROBUTADIENE 1000 U C460 HEXACHLOROBUTADIENE 1000 U C465 P-CHLORO-M-CRESOL 1000 U C465 P-CHLORO-M-CRESOL 1000 U C515 2, 4, 6-TRICHLOROPHENOL 1000 U C515 2, 4, 5-TRICHLOROPHENOL 1000 U C520 2, 4, 5-TRICHLOROPHENOL 5000 U C525 2-CHLORONAPHTHALENE 1000 U C535 DIMETHYL PHTHALATE 1000 U C540 ACENAPHTHYLENE 1000 U C553 3-NITROANILINE 5000 U C555 3-NITROANILINE 5000 U C555 2, 4-DINITROPHENOL 5000 U C555 2, 4-DINITROPHENOL 5000 U C555 2, 4-DINITROPHENOL 5000 U C555 DIBENZOFURAN 1000 U C5570 2, 4-DINITROTOLUENE 1000 U	C370	N-NITROSODIPROPYLAMINE		
C415 ISOPHORONE C420 2-NITROPHENOL C425 2, 4-DIMETHYLPHENOL C430 BENZOIC ACID C430 BIS(2-CHLOROETHOXY)METHANE C440 2, 4-DICHLOROPHENOL C435 BIS(2-CHLOROETHOXY)METHANE C440 1, 2, 4-TRICHLOROBENZENE C440 1, 2, 4-TRICHLOROBENZENE C450 NAPHTHALENE C450 NAPHTHALENE C450 NAPHTHALENE C450 HEXACHLOROBUTADIENE C460 HEXACHLOROBUTADIENE C465 P-CHLORO-M-CRESOL C470 2-METHYLNAPHTHALENE C510 HEXACHLOROCYCLOPENTADIENE C510 HEXACHLOROCYCLOPENTADIENE C510 HEXACHLOROCYCLOPENTADIENE C520 2, 4, 5-TRICHLOROPHENOL C520 2, 4, 5-TRICHLOROPHENOL C530 2-NITROANILINE C530 2-NITROANILINE C530 2-NITROANILINE C540 ACENAPHTHYLENE C540 ACENAPHTHYLENE C550 ACENAPHTHENE C550 ACENAPHTHENE C550 ACENAPHTHENE C550 ACENAPHTHENE C550 ACENAPHTHENE C550 ACENAPHTHENE C550 LIBENZOFURAN C5570 2, 4-DINITROTOLUENE C570 2, 4-DINITROTOLUENE	C375	HEXACHLOROETHANE	1000 U	
C420 2-NITROPHENOL 1000 U C425 2,4-DIMETHYLPHENOL 1000 U C430 BENZOIC ACID 5000 U C435 BIS(2-CHLOROETHOXY)METHANE 1000 U C440 2,4-DICHLOROPHENOL 1000 U C445 1,2,4-TRICHLOROBENZENE 1000 U C445 1,2,4-TRICHLOROBENZENE 1000 U C450 NAPHTHALENE 1000 U C455 4-CHLOROANILINE 1000 U C466 HEXACHLOROBUTADIENE 1000 U C465 P-CHLORO-M-CRESOL 1000 U C470 2-METHYLNAPHTHALENE 1000 U C510 HEXACHLOROCYCLOPENTADIENE 1000 U C515 2,4,6-TRICHLOROPHENOL 1000 U C525 2-CHLOROAPHTHALENE 1000 U C525 2-CHLORONAPHTHALENE 1000 U C530 2-NITROANILINE 5000 U C535 DIMETHYL PHTHALATE 1000 U C540 ACENAPHTHYLENE 1000 U C5545 3-NITROANILINE 5000 U C5555 2,4-DINITROPHENOL 5000 U C5555 2,4-DINITROPHENOL 5000 U C5555 2,4-DINITROPHENOL 5000 U C5565 DIBENZOFURAN 1000 U C5670 2,4-DINITROTOLUENE 1000 U	C410	NI TROBENZENE		
C425 2,4-DIMETHYLPHENOL 1000 U C430 BENZOIC ACID 5000 U C435 BIS(2-CHLORQETHOXY)METHANE 1000 U C440 2,4-DICHLOROPHENOL 1000 U C445 1,2,4-TRICHLOROBENZENE 1000 U C450 NAPHTHALENE 1000 U C455 4-CHLOROBUTADIENE 1000 U C460 HEXACHLOROBUTADIENE 1000 U C465 P-CHLORO-M-CRESOL 1000 U C470 2-METHYLNAPHTHALENE 1000 U C510 HEXACHLOROCYCLOPENTADIENE 1000 U C512 2,4,6-TRICHLOROPHENOL 1000 U C522 2,4,5-TRICHLOROPHENOL 5000 U C533 2-NITROANILINE 5000 U C533 2-NITROANILINE 1000 U C540 ACENAPHTHYLENE 1000 U C551 ACENAPHTHENE 1000 U C552 2,4-DINITROPHENOL 5000 U C553 2,4-DINITROPHENOL 5000 U C555 2,4-DINITROPHENOL 5000 U C555 2,4-DINITROPHENOL 5000 U C556 DIBENZOFURAN 1000 U C565 DIBENZOFURAN 1000 U C5670 2,4-DINITROTOLUENE 1000 U	C415	ISOPHORONE		
C430 BENZDIC ACID  C435 BIS(2-CHLOROETHOXY)METHANE  C440 2, 4-DICHLOROPHENOL  C451 1, 2, 4-TRICHLOROBENZENE  C450 NAPHTHALENE  C450 NAPHTHALENE  C460 HEXACHLOROBUTADIENE  C465 P-CHLORO-M-CRESOL  C470 2-METHYLNAPHTHALENE  C510 HEXACHLOROCYCLOPENTADIENE  C510 HEXACHLOROCYCLOPENTADIENE  C510 2, 4, 6-TRICHLOROPHENOL  C520 2, 4, 5-TRICHLOROPHENOL  C520 2 - NITROANILINE  C530 2-NITROANILINE  C530 2-NITROANILINE  C540 ACENAPHTHYLENE  C550 ACENAPHTHENE  C550 ACENAPHTHENE  C550 ACENAPHTHENOL  C550 ACENAPHTHENOL  C550 ACENAPHTHENOL  C550 ACENAPHTHENOL  C565 DIBENZOFURAN  C570 2, 4-DINITROTOLUENE  C570 C570 C570 C570 C570 C770 C770 C770	C420	2-NITROPHENOL	1000 U	
C430 BENZDIC ACID  C435 BIS(2-CHLOROETHOXY)METHANE  C440 2, 4-DICHLOROPHENOL  C451 1, 2, 4-TRICHLOROBENZENE  C450 NAPHTHALENE  C450 NAPHTHALENE  C460 HEXACHLOROBUTADIENE  C465 P-CHLORO-M-CRESOL  C470 2-METHYLNAPHTHALENE  C510 HEXACHLOROCYCLOPENTADIENE  C510 HEXACHLOROCYCLOPENTADIENE  C510 2, 4, 6-TRICHLOROPHENOL  C520 2, 4, 5-TRICHLOROPHENOL  C520 2 - NITROANILINE  C530 2-NITROANILINE  C530 2-NITROANILINE  C540 ACENAPHTHYLENE  C550 ACENAPHTHENE  C550 ACENAPHTHENE  C550 ACENAPHTHENOL  C550 ACENAPHTHENOL  C550 ACENAPHTHENOL  C550 ACENAPHTHENOL  C565 DIBENZOFURAN  C570 2, 4-DINITROTOLUENE  C570 C570 C570 C570 C570 C770 C770 C770	C425	2,4-DIMETHYLPHENOL	1000 U	
C440 2, 4-DICHLOROPHENOL 1000 U C445 1, 2, 4-TRICHLOROBENZENE 1000 U C450 NAPHTHALENE 1000 U C455 4-CHLOROANILINE 1000 U C460 HEXACHLOROBUTADIENE 1000 U C465 P-CHLORO-M-CRESOL 1000 U C470 2-METHYLNAPHTHALENE 1000 U C510 HEXACHLOROCYCLOPENTADIENE 1000 U C515 2, 4, 6-TRICHLOROPHENOL 1000 U C520 2, 4, 5-TRICHLOROPHENOL 5000 U C525 2-CHLORONAPHTHALENE 1000 U C530 2-NITROANILINE 5000 U C530 DIMETHYL PHTHALATE 1000 U C540 ACENAPHTHYLENE 1000 U C545 3-NITROANILINE 5000 U C550 ACENAPHTHENE 1000 T740 C C550 ACENAPHTHENE 1000 U C560 4-NITROPHENOL 5000 U C560 4-NITROPHENOL 5000 U C565 DIBENZOFURAN 1000 230 C C570 2, 4-DINITROTOLUENE 1000 U	C430		5000 U	
C445 1.2,4-TRICHLOROBENZENE 1000 U C450 NAPHTHALENE 1000 U C455 4-CHLOROANILINE 1000 U C460 HEXACHLOROBUTADIENE 1000 U C465 P-CHLORO-M-CRESOL 1000 U C470 2-METHYLNAPHTHALENE 1000 U C510 HEXACHLOROCYCLOPENTADIENE 1000 U C511 2,4,6-TRICHLOROPHENOL 1000 U C520 2,4,5-TRICHLOROPHENOL 5000 U C525 2-CHLORONAPHTHALENE 1000 U C530 2-NITROANILINE 5000 U C530 DIMETHYL PHTHALATE 1000 U C540 ACENAPHTHYLENE 1000 . 170 C C545 3-NITROANILINE 5000 U C550 ACENAPHTHENE 1000 . 740 C C550 2,4-DINITROPHENOL 5000 U C560 4-NITROPHENOL 5000 U C565 DIBENZOFURAN 1000 . 230 C C570 2,4-DINITROTOLUENE 1000 U	C435			
C450 NAPHTHALENE C455 4-CHLQRQANILINE C460 HEXACHLQROBUTADIENE C465 P-CHLQRO-M-CRESQL C470 2-METHYLNAPHTHALENE C510 HEXACHLQRQCYCLQPENTADIENE C510 HEXACHLQRQCYCLQPENTADIENE C515 2, 4, 6-TRICHLQRQPHENQL C520 2, 4, 5-TRICHLQRQPHENQL C520 2, 4, 5-TRICHLQRQPHENQL C530 2-NITRQANILINE C530 2-NITRQANILINE C540 ACENAPHTHYLENE C540 ACENAPHTHYLENE C550 ACENAPHTHENE C550 ACENAPHTHENE C550 ACENAPHTHENE C550 ACENAPHTHENE C550 4-NITRQPHENQL C560 4-NITRQPHENQL C565 DIBENZQFURAN C570 2, 4-DINITRQTQLUENE C570 2, 4-DINITRQTQLUENE C570 2, 4-DINITRQTQLUENE C570 C570 2, 4-DINITRQTQLUENE C570 C570 C570 C570 C770 C770 C770 C770	C440	2,4-DICHLOROPHENOL	1000 U	
C455 4-CHLOROBATLINE 1000 U C460 HEXACHLOROBUTADIENE 1000 U C465 P-CHLORO-M-CRESOL 1000 U C470 2-METHYLNAPHTHALENE 1000 . 1500 C510 HEXACHLOROCYCLOPENTADIENE 1000 U C515 2, 4, 6-TRICHLOROPHENOL 1000 U C520 2, 4, 5-TRICHLOROPHENOL 5000 U C525 2-CHLORONAPHTHALENE 1000 U C530 2-NITROANILINE 5000 U C535 DIMETHYL PHTHALATE 1000 U C540 ACENAPHTHYLENE 1000 . 170 C C545 3-NITROANILINE 5000 U C555 2, 4-DINITROPHENOL 5000 U C556 4-NITROPHENOL 5000 U C566 10BENZOFURAN 1000 . 230 C C570 2, 4-DINITROTOLUENE 1000 U		1.2.4-TRICHLOROBENZENE	1000 U	
C460 HEXACHLOROBUTADIENE 1000 U C465 P-CHLORO-M-CRESOL 1000 U C470 2-METHYLNAPHTHALENE 1000 1500 C510 HEXACHLOROCYCLOPENTADIENE 1000 U C515 2, 4, 6-TRICHLOROPHENOL 1000 U C520 2, 4, 5-TRICHLOROPHENOL 5000 U C525 2-CHLORONAPHTHALENE 1000 U C530 2-NITROANILINE 5000 U C535 DIMETHYL PHTHALATE 1000 U C540 ACENAPHTHYLENE 1000 170 C C545 3-NITROANILINE 5000 U C550 ACENAPHTHENE 1000 740 C C555 2, 4-DINITROPHENOL 5000 U C566 DIBENZOFURAN 1000 230 C C570 2, 4-DINITROTOLUENE 1000 U			1000	. 5100
C445 P-CHLORO-M-CRESOL  C470 2-METHYLNAPHTHALENE  C510 HEXACHLOROCYCLOPENTADIENE  C515 2,4,6-TRICHLOROPHENOL  C520 2,4,5-TRICHLOROPHENOL  C525 2-CHLORONAPHTHALENE  C530 2-NITROANILINE  C540 ACENAPHTHYLENE  C540 ACENAPHTHYLENE  C545 3-NITROANILINE  C550 ACENAPHTHENE  C550 ACENAPHTHENE  C550 ACENAPHTHENE  C550 ACENAPHTHENE  C550 4-DINITROPHENOL  C560 4-NITROPHENOL  C565 DIBENZOFURAN  C570 2,4-DINITROTOLUENE  1000			1000 U	
C470 2-METHYLNAPHTHALENE 1000			1000 U	
C510 HEXACHLOROCYCLOPENTADIENE 1000 U C515 2, 4, 6-TRICHLOROPHENOL 1000 U C520 2, 4, 5-TRICHLOROPHENOL 5000 U C525 2-CHLORONAPHTHALENE 1000 U C530 2-NITROANILINE 5000 U C535 DIMETHYL PHTHALATE 1000 U C540 ACENAPHTHYLENE 1000 170 C C545 3-NITROANILINE 5000 U C550 ACENAPHTHENE 1000 740 C C555 2, 4-DINITROPHENOL 5000 U C560 4-NITROPHENOL 5000 U C565 DIBENZOFURAN 1000 230 C C570 2, 4-DINITROTOLUENE			1000 U	
C515 2,4,6-TRICHLOROPHENOL 1000 U C520 2,4,5-TRICHLOROPHENOL 5000 U C525 2-CHLORONAPHTHALENE 1000 U C530 2-NITROANILINE 5000 U C535 DIMETHYL PHTHALATE 1000 U C540 ACENAPHTHYLENE 1000 . 170 C C545 3-NITROANILINE 5000 U C550 ACENAPHTHENE 1000 . 740 C C555 2,4-DINITROPHENOL 5000 U C565 DIBENZOFURAN 1000 . 230 C C570 2,4-DINITROTOLUENE 1000 U			1000	. 1500
C520 2,4,5-TRICHLOROPHENOL 5000 U C525 2-CHLORONAPHTHALENE 1000 U C530 2-NITROANILINE 5000 U C535 DIMETHYL PHTHALATE 1000 U C540 ACENAPHTHYLENE 1000		· · · · · · · · · · · · · · · · · · ·		
C525 2-CHLORONAPHTHALENE C530 2-NITROANILINE C535 DIMETHYL PHTHALATE C540 ACENAPHTHYLENE C545 3-NITROANILINE C550 ACENAPHTHENE C550 ACENAPHTHENE C555 2, 4-DINITROPHENOL C560 4-NITROPHENOL C565 DIBENZOFURAN C570 2, 4-DINITROTOLUENE 1000 U C560 U C570 2, 4-DINITROTOLUENE				
C530 2-NITROANILINE 5000 U C535 DIMETHYL PHTHALATE 1000 U C540 ACENAPHTHYLENE 1000				
C535 DIMETHYL PHTHALATE 1000 U C540 ACENAPHTHYLENE 1000				
C540 ACENAPHTHYLENE 1000				
C545 3-NITROANILINE 5000 U C550 ACENAPHTHENE 1000 740 C C555 2,4-DINITROPHENOL 5000 U C560 4-NITROPHENOL 5000 U C565 DIBENZOFURAN 1000 230 C C570 2,4-DINITROTOLUENE 1000 U				
C550 ACENAPHTHENE 1000 740 C C555 2,4-DINITROPHENOL 5000 U C560 4-NITROPHENOL 5000 U C565 DIBENZOFURAN 1000 230 C C570 2,4-DINITROTOLUENE 1000 U		** =		. 170 J
C555 2,4-DINITROPHENOL 5000 U C560 4-NITROPHENOL 5000 U C565 DIBENZOFURAN 1000 230 C C570 2,4-DINITROTOLUENE 1000 U		<del>-</del>		
C560 4-NITROPHENOL 5000 U C565 DIBENZOFURAN 1000 230 C C570 2,4-DINITROTOLUENE 1000 U				. 740 J
C565 DIBENZOFURAN 1000 230 C C570 2,4-DINITROTOLUENE 1000 U				
C570 2,4-DINITROTOLUENE 1000 U				
		<del>-</del>		. 230 J
C543 2,6-DINITROTOLUENE 1000 U				
	C543	2,6-DINITROTOLUENE	1000 U	

SAMPLE NUMBER: SPECO001-01 SI-16

. SPECOOOI-O1 31 18

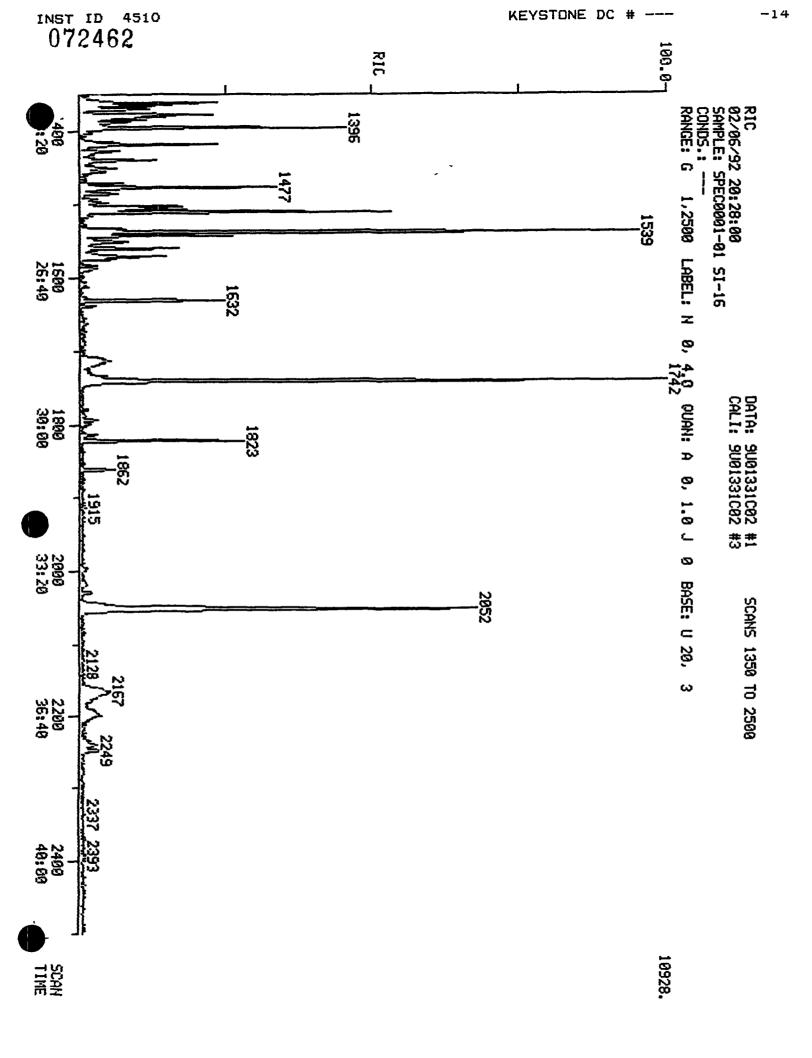
VOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 9001331C02

 - <del></del>	COMPOUND	DETECTION LIMIT	AMOUNT FOUND
 		(MILLIGRAMS / I	KG) 
C580	DIETHYL PHTHALATE	1000 U	
C585	4-CHLOROPHENYL PHENYL ETHER		
C590	FLUORENE	1000	1400
C595	4-NITROANILINE	5000 U	1400
C610	4,6-DINITRO-2-METHYLPHENOL	5000 U	
C615	N-NITROSODIPHENYLAMINE	1000 U	
C625	4-BROMOPHENYL PHENYL ETHER		
C630	HEXACHLOROBENZENE	1000 U	
C635	PENTACHLOROPHENOL	5000 U	
C640	PHENANTHRENE	1000	2700
C645	ANTHRACENE	1000	250 J
C650	DI-N-BUTYL PHTHALATE	1000 U	
C655	FLUORANTHENE	1000	400 J
C715	PYRENE	1000	720 J
C720	BUTYL BENZYL PHTHALATE	1000 U	
C725	3,3'-DICHLOROBENZIDINE	2000 U	
C730	BENZO(A)ANTHRACENE	1000 U	
C745	BIS(2-ETHYLHEXYL)PHTHALATE	1000 U	
C740	CHRYSENE	1000 U	
C760	DI-N-OCTYL PHTHALATE	1000 U	
C765	BENZO(B)FLUORANTHENE	1000 U	
C770	BENZO(K)FLUORANTHENE	1000 U	
C775	BENZO(A)PYRENE	1000 U	
C780	INDENO(1, 2, 3-CD)PYRENE	1000 U	
C785	DIBENZO(A, H) ANTHRACENE	1000 U	
C790	BENZO(GHI)PERYLENE	1000 U	

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT



SACT 2463 SPECO001-01 SI-16

FILENAME 9001331C02 INST ID 4510

CLIENT FRENCH ANALYST JDD

E INJECTED 02/06/92 20:28 00 VERIFIED BY

CALIB STD DATE 01/27/92 CORR. FACTOR 100 00

SCAN#		SEMIVOLATILE COMPOUNDS	M/E	AMOUNT	RRT	AREA
495	CI30	1,4-DICHLOROBENZENE-D4 ** IS	152	2000 MG/KG	1.000	10729
727	C140	NAPHTHALENE-D8 ** IS2 **	136	2000 MG/KG	1.000	46761
1028	C150	ACENAPHTHENE-D10 ** IS3 **	164	2000 MG/KG	1.000	23665
1275	C160	PHENANTHRENE-D10 ** IS4 **	188	2000 MG/KG	1.000	26162
1742	C170	CHRYSENE-D12 ** IS5 **	240	2000 MG/KG	1.000	16098
2052	C175	PERYLENE-D12 ** IS6 **	264	2000 MG/KG	1.000	13207
730	C450	NAPHTHALENE	128	5040 MG/KG	1.004	131380
853	C470	2-METHYLNAPHTHALENE	142	1504 MG/KG	1. 173	24400
1000	C540	ACENAPHTHYLENE	152	169 MG/KG	0. 973	3546
1032	C550	ACENAPHTHENE	153	741 MG/KG	1.004	9753
1061	C545	DIBENZOFURAN	168	227 MG/KG	1.032	4077
1121	C590	FLUORENE	166	1414 MG/KG	1.090	20034
1279	C640	PHENANTHRENE	178	2710 MG/KG	1.003	46050
1285	C645	ANTHRACENE	178	248 MG/KG	1.008	3527
1477	C655	FLUORANTHENE	202	404 MG/KG	1.158	5564
1511	C715	PYRENE	202	717 MG/KG	0. 867	8492

#### ORGANICS ANALYSIS DATA SHEET -

DRATORY NAME: KEYSTONE

SAMPLE ID NO.: 91---

SAMPLE MATRIX: SOIL

DATA RELEASE AUTHORIZED BY.

CASE NO.: ---

QC REPORT NO.: 01/27/92

CONTRACT NO. . ---

DATE SAMPLE RECEIVED: ... . .....

SEMIVOLATILES

CONCENTRATION: LOW DATE EXTRACTED: ......

DATE ANALYZED: 02/07/92

DATAFILE: 9B0131C01W

	COMPOUND	DETECTION LIMIT (MILLIGRAM	FOUND
C315		100 U	، نشبیه نید موجه کا که نیا نید ها
C325	BIS(2-CHLOROETHYL)ETHER	100 U	
C330		100 U	
C335	1,3-DICHLOROBENZENE	100 U	
C340	1,4-DICHLOROBENZENE	100 U	
C345	BENZYL ALCOHOL	100 U	
C350	1,2-DICHLOROBENZENE	100 U	
C355	1.4-DICHLOROBENZENE BENZYL ALCOHOL 1.2-DICHLOROBENZENE 2-METHYLPHENOL	100 U	
C357	2,2'-OXYBIS(1-CHLOROPROPANE)	100 U	
C365	4-METHYLPHENOL	100 U	
C370	N-NITROSODIPROPYLAMINE	100 U	
C375	HEXACHLOROETHANE	100 U	
C410	NITROBENZENE	100 U	
C415	ISOPHORONE	100 U	
C420	2-NITROPHENOL	100 U	
C425	2,4-DIMETHYLPHENOL BENZOIC ACID	100 U	
C430	BENZOIC ACID	500 U	
C435	BIS(2-CHLOROETHOXY)METHANE	100 U	
C440	2.4-DICHLOROPHENOL	100 U	
C445		100 U	
C450		100 U	
C455		100 U	
C460		100 U	
C465		100 U	
C470		100 U	
C510		100 U	
C515	· ··· · · · · · · · · · · · · · · · ·	100 U	
C520		500 U	
C525		100 U	
C530		500 U	
C535		100 U	
	HOLIANI HILLI ELIAL	100 U	
C545	3-NITROANILINE	500 U	
C550		100 U	
C555		500 U	
C560	4-NITROPHENOL	500 U	
C565	DIBENZOFURAN	100 U	
C570	= =	100 U	
C543	2,6-DINITROTOLUENE	100 U	

SAMPLE NUMBER . SBLK WD

100 U

-----

..IVOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

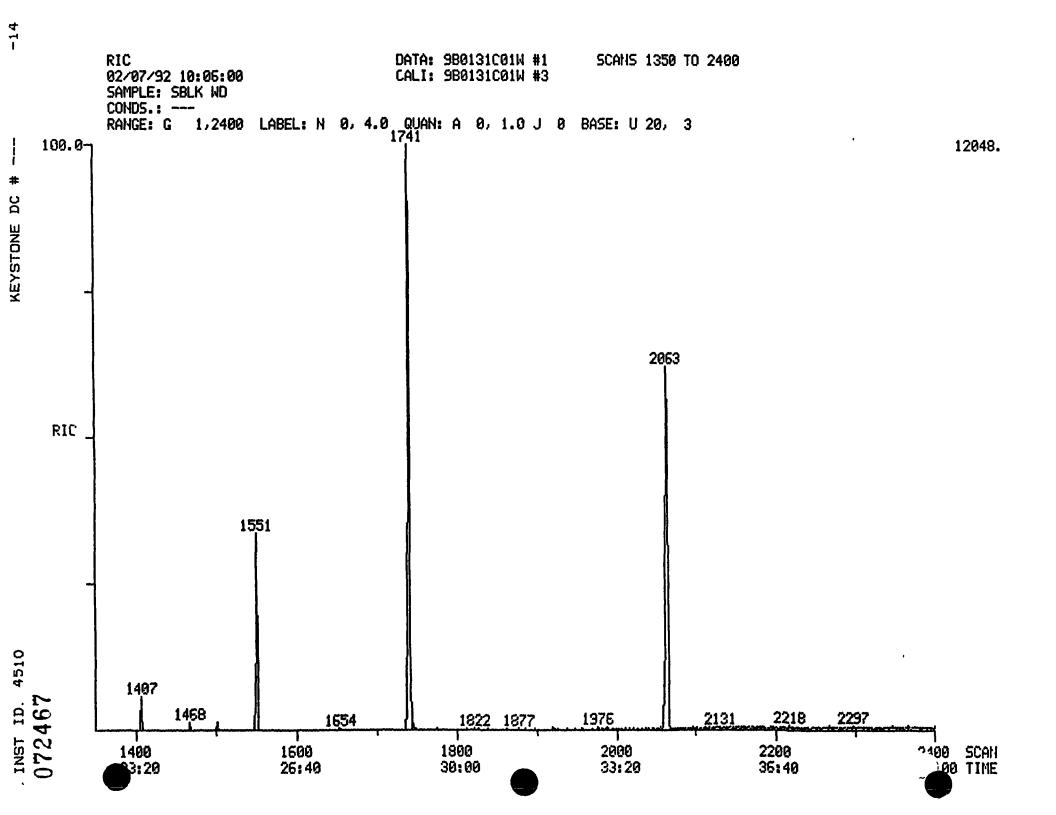
DATAFILE: 9B0131C01W

		COMPOUND	DETECTION LIMIT (MILLIGRAMS	AMOUNT FOUND / KG)
	C580	DIETHYL PHTHALATE	100 U	
	C585		100 U	
	C590		100 U	
	C595	4-NITROANILINE	500 U	
	C610	4,6-DINITRO-2-METHYLPHENOL	500 U	
	C615	N-NITROSODIPHENYLAMINE	100 U	
	C625	4-BROMOPHENYL PHENYL ETHER	100 U	
	C630	HEXACHLOROBENZENE	100 U	
	C635	PENTACHLOROPHENOL	500 U	
	C640	PHENANTHRENE	100 U	
	C645	ANTHRACENE	100 U	
	C&50	DI-N-BUTYL PHTHALATE	100 U	
	C655	FLUORANTHENE	100 U	
	C715	PYRENE	100 U	
_	C720	BUTYL BENZYL PHTHALATE	100 U	
	C725	3,3'-DICHLOROBENZIDINE	200 U	
	C730	BENZO(A)ANTHRACENE	100 U	
	C745	BIS(2-ETHYLHEXYL)PHTHALATE	100 U	
	C740	CHRYSENE	100 U	
	<b>C760</b>	DI-N-OCTYL PHTHALATE	100 U	
	C765	BENZO(B)FLUORANTHENE	100 U	
	C770	BENZO(K)FLUORANTHENE	100 U	
	C775	BENZO(A)PYRENE	100 U	
	C780	INDENO(1, 2, 3-CD)PYRENE	100 U	

U = UNDETECTED AT THE LISTED DETECTION LIMIT

C785 D(BENZO(A, H)ANTHRACENE C790 BENZO(GHI)PERYLENE

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT



5077254 AB

SBLK WD

FILENAME

9B0131C01W

INST ID

4510

CLIENT

KEYSTONE

ANALYST

JDD

E INJECTED

02/07/92 10:06:00

VERIFIED BY

CALIB STD DATE

01/27/92

CORR. FACTOR

10.00

SCAN#		SEMIVOLATILE COMPOUNDS	 M/E	AMOUNT	RRT	AREA
30AN#	· ·	SENIVULATILE CONFOUNDS	177E			
492	C130	1,4-DICHLOROBENZENE-D4 ** IS	152	200 MG/KG	1.000	9358
722	CI40	NAPHTHALENE-DB ** IS2 **	136	200 MG/KG	1.000	37050
1021	CI50	ACENAPHTHENE-D10 ** IS3 **	164	200 MG/KG	1.000	16602
1267	C140	PHENANTHRENE-D10 ** IS4 **	188	200 MG/KG	1.000	19525
1741	C170	CHRYSENE-D12 ** IS5 **	240	200 MG/KG	1.000	14648
2063	CI75	PERYLENE-D12 ** IS6 **	264	200 MG/KG	1.000	12206
278	<b>CS50</b>	2~FLUOROPHENOL ** SU1 **	112	75 15% 000	0. 565	6728
507	<b>CS45</b>	PHENOL-D5 ** SU2 **	99	65 23 7	1.030	8795
1160	CS55	2,4,6-TRIBROMOPHENOL ** SU5	330	40 8%	1, 136	200
476	<b>CS70</b>	2-CHLOROPHENOL-D4 ** SU7 **	132	70 14%	0. 967	8324
526	<b>CS75</b>	1,2-DICHLOROBENZENE-D4 ** SU	150	80 15%	1.069	5325
600	CS20	NITROBENZENE-D5 ** SU3 **	82	85 17%	0. 831	5234
920	CS25	2-FLUOROBIPHENYL ** SU4 **	172	95 19%	0. 901	8755
1551	CS30	TERPHENYL-D14 ** SU6 **	244	70 44 %	0. 891	5517

No.: 86 Page 1 of 2

REPORTING LABORATORY FRENCH LTD. PROJECT Name: Keystone Lab-Houston LIG, Incorporated 15010 FM 2100, Suite 200 Crosby, Texas 77532 Address: 8300 Westpark Houston, TX 77063 Contact: Daniel D. Pastalaniec Phone: (713) 266-6800 FAX: (713) 974-5494 (713) 328-5860 FAX (713) 328-2996 Requested By: TD DATA PACKAGE TO: FLTG, Incorporated Crosby, TX 77532 1024 Gulf Pump Road Days: Standard TA: n \*\* Notify Lab Contact Immediately \*\* Date: Notified by: FLTG MATRIX CODE: SPEC FLTG SET NUMBER: SPECO001 Quality Control Level: I Site Location: 1024 Gulf Pump Road Crosby, TX 77532 Grab/ FLTG Location Matrix Time Sample No Date Comp Type 08:45 SI-16 Grab Water **ENV** SPEC0001-01 1/30/92 Ship in separate container SAMPLER'S Name(s) (PRINT) Affiliation 1. ROKET OLIVER Operator GAF 3. Carrier: Bill No.: RELINQUISHED BY RECEIVED BY (SIGN) (INITIALS) Date Time (SIGN) (INITIALS) Date Time 1. Robert Oliver RD. 1-30-92 3:27 1/35/91 3.2) RETURNED BY DISPOSED BY (INITIALS) Date Time Time (SIGN) (SIGN) (INITIALS) Date 2. NOTES TO LAB: Sample to be analyzed per DAF

# Analysis Request and Chain Custody Record

No.: 86 Page 2 of 2

LTG MATRIX (	"LTG MATRIX CODE:SPEC Lab Work Order No:					
SAMPLING SOP	REFERENCE:					~ ~ ~ ~
DESCRIPTION:	DESCRIPTION: Groundwater					
NOTES TO SAMI	PLER:					
	P	ARAMETER	S REQUEST	ED		
voa svoa tox teh						
FLTG Sample No	Lab Sample ID	Conta	iner		Preserv	ative
SPEC0001-01	<u></u>	2-4 oz	bottle		refrig	
				{ 		
		 	·			
COMMENTS			Set (SIGN)	QC Complet (INIT	ed by: TALS)	Date

# BOOKNYKK

# Appendix 2

Analytical Reports INT-11 DNAPL sample

01502434

Report Prepared for

FLTG, INC.

1024 GULF PUMP ROAD. CROSBY TX 77532

Attention : TED DAVIS

by

Keystone Lab - Houston
A Division of CHESTER LabNet
8300 Westpark Drive
Houston, Texas 77063
(713) 266-6800

CERTIFIED BY :

Dan Pastalaniec Project Manager

> WORK ORDER : H92-04.108 DATE RECEIVED : 9-APR-1992

PROJECT ID : P.O. NUMBER : SPEC0015

### SAMPLE & ANALYSIS SUMMARY

Client's Sample Number			Sample Matrix
LAB BLANK SPEC0015-01 SPEC0015-01DUP SPEC0015-01MS	04/08/92	14:40	WATER WATER WATER WATER
V\$TCL	Time		
	<del></del>		<del> </del>
	Time		
Client	Client	NM::1+in1	ier
		-T	
· · · · · · · · · · · · · · · · · · ·		· 	<del></del>
· · · · · · · · · · · · · · · · · · ·	×	'. <del></del>	
		1 	<del></del>
TENT			
ID   Wt. %			
	Sample Number  LAB BLANK SPEC0015-01 SPEC0015-01MS  06/18/92 04/09/92 10:35  Report V\$TCL Analyst   Date-  Report CB Analyst   Date-  I Multipliers Client   Sample ID Multipliers   -01   1	LAB BLANK SPEC0015-01 04/08/92 SPEC0015-01DUP 04/08/92 SPEC0015-01MS 04/08/92 06/18/92 FLTG Set Number 04/09/92 10:35 Lab Number: HS  Report V\$TCL  Analyst   Date-Time  Report CB  Analyst   Date-Time  Client Client   Sample ID   Multiplier   Sample ID     -01	Sample Number   Collected

COMMENTS:

#### SET NARRATIVE Matrix SPEC

Keystone Lab-Houston

FLTG

Lab Work Order: H92-04.108

Project Set Number: SPEC0015

Report Submission Date: 06/18/92

I. SAMPLE COLLECTION/SHIPPING PROBLEMS

FLTG

Lab

Sample No. Param. Sample No. Problem

Outcome

NONE

II. ANALYSIS PROBLEMS

FLTG

Lab

Sample No. Param.

Sample No. Problem

Outcome

NONE

III. QUALITY ASSURANCE PROBLEMS

FLTG

Sample No.

File No. Problem

Outcome

SPEC0015-01

4U04108V02 SU 1 Not Found ; SU 2 >110% Matrix Effec

4U04108V02 SU 1 >620% Matrix Effect

SPEC0015-01

COMMENTS- MS/MSD on VOA not done because of Matrix. Sample was run once within holding time but the data was worthless due to the high contamination. The three submitted runs were one day after the hold time expired.

Page 1

17-JUN-1992

## Summary of Analytical Results

Date received: 9-APR-1992

Oustoner: FLIG , INC.

Job name: H92-04.108

		Samples			•	
Keystone ID Sampling Point Date Sampled Customer ID		108-001 CA CC 8-APR-1992 IAB BLANK	108-002 X 8-APR-1992 SPEC0015-01 NA	108-003 QA QC 8-APR-1992 SPEC0015-01 DUP	108-004 QA_QC 8-APR-1992 SPECCO15-01 MS	
Parameters	Units			<del> </del>		
Total Organic Carbon Analyst: RNS Date/Time: 04/13/92 10 Dilution: 10	mg/L ):31	₫.0	506	556	97.0*	
Total Organic Halogens Analyst: DH	mg/L	<0.005	595	599	103*	

Date/Time: 04/14/92 14:00 Dilution: 10000

\* - % Recovery NR - Not Required NA - Not Applicable

# SAMPLE NUMBER: SPECO015-01 INT 11 1

#### DRGANICS ANALYSIS DATA SHEET

ARATORY NAME: KEYSTONE ENV. AB SAMPLE ID NO.: 910410802

CASE NO. : ----INITIAL CALIBRATION DATE: 04/22/92

AMPLE MATRIX: WATER

CONTRACT NO.: ---ATA RELEASE AUTHORIZED BY:..... DATE SAMPLE RECEIVED: 04/09/92

#### VOLATILES

DATAFILE: 4U04108V02B CONCENTRATION: LOW ATE ANALYZED: 04/23/92 DILUTION FACTOR: 10000.00

	, <del></del>		DETECTION AMOUNT LIMIT FOUND (MICROGRAMS / LITER )	
	C010	CHLOROMETHANE	100000 U	
	CO15	BROMOMETHANE	100000 U	
	CO20	BROMOMETHANE VINYL CHLORIDE	100000 U	
	CO25	CHLOROETHANE METHYLENE CHLORIDE	100000 U	
	CO30	METHYLENE CHLORIDE	50000 20000 J	
	C035	ACETONE	100000 U	
	C040	CARBON DISULFIDE	50000 U	
	C045	1,1-DICHLOROETHENE	50000 U 50000 U 50000 20000 J	
	C050	1,1-DICHLOROETHANE	50000 20000 J 50000 21000 J 50000 810000	
	C053	1,1-DICHLOROETHANE 1,2-DICHLOROETHENE (TOTAL)	50000 21000 J	
V	C040	CHLOROFORM		
	C065	1,2-DICHLORGETHANE	50000650000	
	C110	2-BUTANONE	100000 23000 J	
	C115	1,1,1-TRICHLORGETHANE		
	C120	CARBON TETRACHLORIDE		
	C125	VINYL ACETATE	100000150000	
	C130	BROMODICHLOROMETHANE	50000 U	
	C140	I'S-DICHFOKOLKOLWE	30000 0	
	C143	CIS-1, 3-DICHLOROPROPENE		
	C150	TRICHLOROETHENE	50000 U	
	C155	DIBROMOCHLOROMETHANE	50000 U	
	C160	1,1,2-TRICHLORDETHANE	50000 U	
	C165	BENZENE	50000 4400 J	
	C172	TRANS-1, 3-DICHLOROPROPENE		
	C180	BROMDFORM	50000 U	
	C205	4-METHYL-2-PENTANONE	100000 27000 J	
	C210	2-HEXANONE TETRACHLOROETHENE 1, 1, 2, 2-TETRACHLOROETHANE	100000 U	
	C550	TETRACHLOROETHENE	50000 18000 J	
	C225	TETRACHLOROETHENE  1, 1, 2, 2-TETRACHLOROETHANE	50000 19000 J	
	C530	TULUENE	20000 0	
	C235	CHLOROBENZENE	50000 U	
	C240	ETHYLBENZENE	50000 U	
	C245	STYRENE	50000 U	
	C250	XYLENES (TOTAL)	50000 U	

UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

NST 7247790

CAS #

ORGANICS ANALYSIS DATA SHEET - PAGE 4

ABORATORY NAME: KEYSTONE ENVIRONMENTAL RESOURCES, INC. CASE NO.: ---
3C REPORT NO.: ..... ANALYST: PVG DATAFILE: 4U04108V02

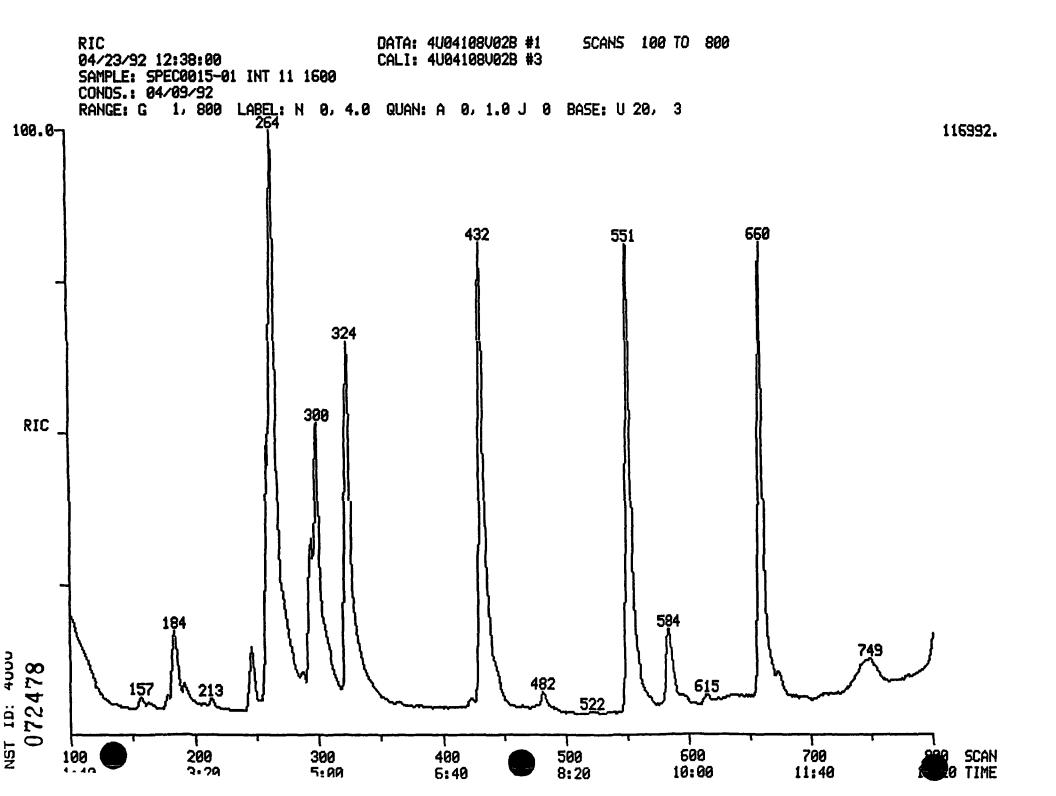
B. TENTATIVELY IDENTIFIED COMPOUNDS

UNKNOWN 184 250000

SCAN# PURITY AMOUNT

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

VOLATILE COMPOUND NAMES



FRENCH LTD. PROJECT REPORTING LABORATORY FLTG, Incorporated Name: Keystone Lab-Houston \$5010 FM 2100. Suite 200 Address: 8300 Westpark rosby, Texas 77532 Houston, TX 77063 Contact: Daniel D. Pastalaniec Phone: (713) 266-6800 (713) 328-5860 FAX (713) 328-2996 (713) 974-5491 FAX: DATA PACKAGE TO: Requested By: TD FLTG. Incorporated Standard TA: Y Days: 21 Crosby, TX 77532 1024 Gulf Pump Road Notified by: Date: FLTG MATRIX CODE: SPEC FLTG SET NUMBER: SPECO015 Quality Control Level: I Site Location: 1024 Gulf Pump Road Crosby, TX 77532 FLTG Grab/ Sample No Date Time Location Comp Matrix Type 1440 SPEC0015-01 4/08/92 **INT 11** Grab Water ENV 7500 SAMPLER'S Name(s) (PRINT)

1. Jeffrey H Barker Affiliation Operator G#F Technical Services 2. 3. Bill No .: Carrier: RECEIVED BY RELINGUISHED BY (SIGN) (INITIALS) (INITIALS) Date (SIGN) Date Time Time 4/9/92 10:35 hun 4/9/42 10:35 DISPOSED BY RETURNED BY (SIGN) (INITIALS) Date Time (SIGN) (INITIALS) Date Time 1. 2. NOTES TO LAB:

FLTG MATRIX CODE:SPEC Lab Work Order No: FLTG SET NUMBER:SPEC0015				
AMPLING SOP	REFERENCE: Wate	r		
DESCRIPTION:	Water			
NOTES TO SAM	PLER:			
	Р	ARAMETER	S REQUESTED	
TOC	т	ox	VOA\$TCL	
FLTG Sample No	Lab Sample ID	Conta	iner	Preservative
SPEC0015-01 SPEC0015-01		2-40ml 4 oz g	lass (z)	HCL H2SO4
SPEC0015-01		8 oz g	lass (1)	HCL
		i i		
COMMENTS			Set QC Complet (SIGN) (INIT	ed by: IALS) Date

# BOOKMARK

Appendix 3

Cone Penetrometer Logs S1-16 Area

01502435

#### **FUGRO GEOSCIENCES, INC.**



6105 Rookin Houston, Texas 77074 Tel· (713) 778-5580 Fax (713) 778-5501

May 18, 1992 Report Number 0301-2126

FLTG, Incorporated 15010 FM 2100, Suite 200 Crosby, Texas 77532

Attention: Mr Mike Day / Mr Bill Edmondson

CONE PENETROMETER TESTING
AND RELATED SERVICES
FOR AFFECTED GROUNDWATER DELINEATION
FRENCH LIMITED SITE
CROSBY, TEXAS

#### Gentlemen

Please find enclosed herewith the final results of the cone penetrometer tests conducted at the above referenced location

For your information, the soil stratigraphy was identified using Campanella and Robertson's Simplified Soil Behavior Chart. Please note that because of the empirical nature of the soil behavior chart, the soil identification should be verified locally

Fugro Geosciences appreciates the opportunity to be of service to your organization if you should have any questions, or if we can be of further assistance, please do not hesitate to contact us. We look forward to working with you in the future.

Very truly yours,

**FUGRO GEOSCIENCES, INC.** 

L David Parker, P E.

**CPT Department Manager** 

LDP/kjp

## **Key To Soil Classification and Symbols**

# SOIL TYPE (Shown in Symbol Column) Sand Silt Clay Fill Sandy Silty Clayey Predominant Type Shown Heavy

#### TERMS DESCRIBING CONSISTENCY OR CONDITION

#### COARSE GRAINED SOILS (Major portion Retained on No 200 Sieve)

includes (1) clean gravels and sand described as fine, medium or course, depending on distribution of grain sizes (2) silty or clayey gravels and sands and (3) fine grained low plasticity soils (PI < 10) such as sandy silts. Condition is rated according to relative density, as determined by lab tests or estimated from resistance to sampler penetration.

Descriptive Term	Penetration Resistance*	Relative Density
Loose	0 - 10	0 to 40%
Medium Dense	10 - 30	40 to 70%
Dense	30 - 50	70 to 90%
Very Dense	Over 50	90 to 100%

<sup>\*</sup> Blows/Foot, 140# Hammer, 30" Drop

#### FINE GRAINED SOILS (Major Portion Passing No 200 Sieve)

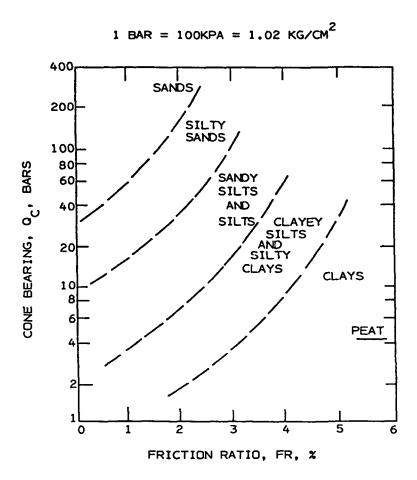
includes (1) inorganic and organic silts and clays, (2) sandy, gravelly or silty clays, and (3) clayey silts. Consistency is rated according to shearing strength, as indicated by penetrometer readings or by unconfined compression tests for soils with PI > 10

Descriptive	Cohesive Shear Strength Tons/Square Foot		
Term_			
Very Soft	Less Than 0 125		
Soft	0 125 to 0 25		
Firm	0 25 to 0 50		
Stuff	0 50 to 1 00		
Very Stiff	1 00 to 2 00		
Hard	2 00 and Higher		

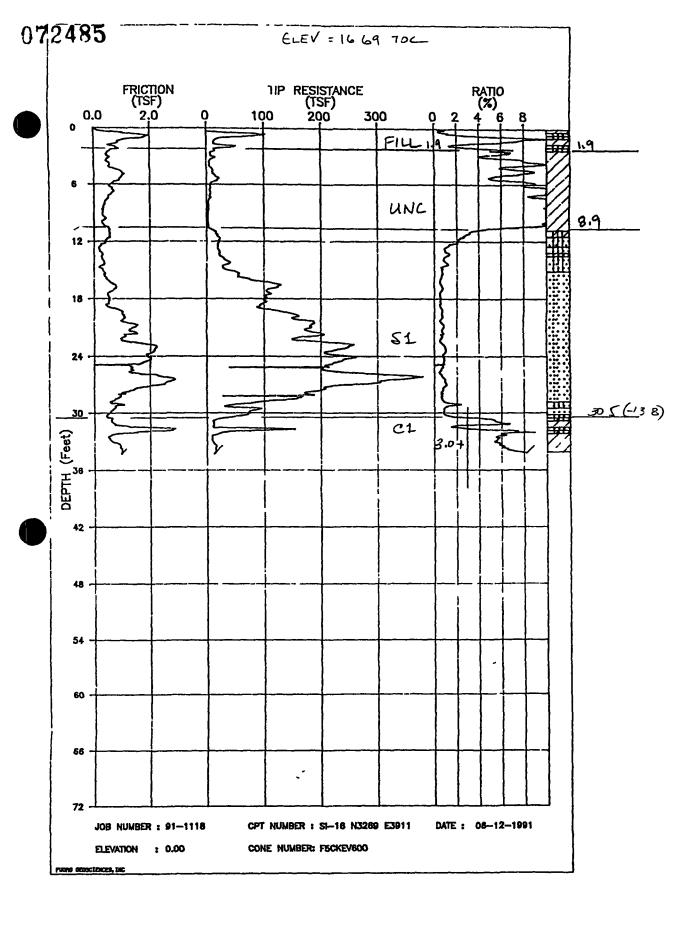
Note Slickensided and fissured clay may have lower unconfined compressive strengths than shown above because of planes of weakness or shrinkage cracks, consistency ratings of such soils are based on hand penetrometer readings

#### TERMS CHARACTERIZING SOIL STRUCTURE

Parting	paper thin in size		pertaining to cohesive soils that exhibit a loose knit or flakey structure
Seam	1/8" to 3" thick		having inclined planes of weakness that are
Layer	greater than 3"		• ,
Fissured	containing shrinkage cracks, frequently filled with	•	slick and glossy in appearance
	fine sand or silt, usually more or less vertical	Degree of Slickensided [	Development
Sensitive	pertaining to cohesive soils that are subject to appreciable loss of strength when remolded	Slightly Slickensided	slickensides present at intervals of 1' to
interbedded	composed of alternate layers of different soil types		2', soil does not easily break along these plates
Laminated	composed of thin layers of varying color and texture	Moderately Slickensided	slickensides spaced at intervals of 1' to 2', soil breaks easily along these planes
Calcareous	containing appreciable quantities of calcium carbonate	Extremely Slickensided.	continuous and interconnected slicken- sides spaced at intervals of 4" to 12',
Well Graded	having wide range in grain sizes and substantial amounts of all intermediate particle sizes		soil breaks along the slickensides into pieces 3" to 6" in size
Poorly Graded	predominantly of one grain size, or having a range of sizes with some intermediate size missing	Intensely Slickensided	slickensides spaced at intervals of less than 4", continuous in all directions, soil breaks down along planes into nodules 1/4" to 2" in size

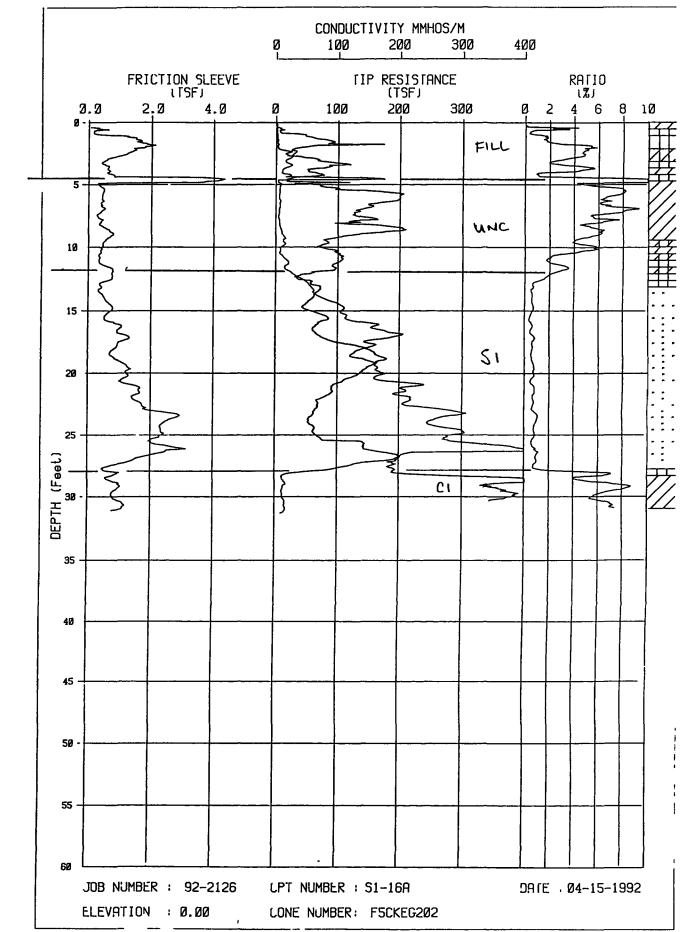


CAMPANELLA AND ROBERTSON CLASSIFICATION CHART

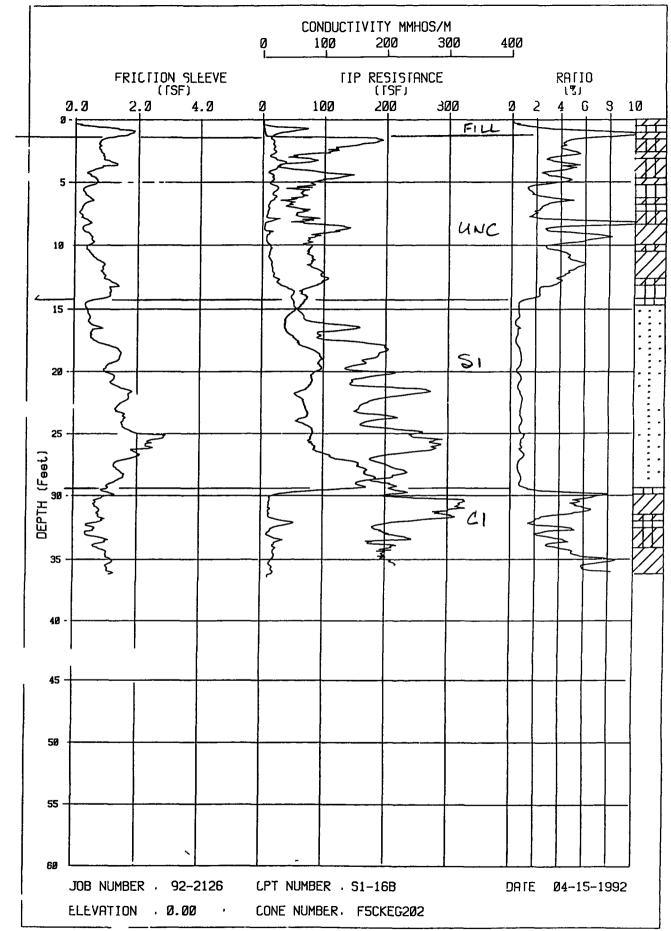


WELL CONSTRUCTION LOC FRENCH, LTD. SUPERFUNDS	CONTR	ACTOR: LAYNE	BECHTEL JOB NUMBER: 21454-100	WELL NUMBER S1-16
GEOLOGIST: DATE	BEGUN: DA	TE COMPLETED:	COORDINATES:	PAGE 1 OF 1
M.A. GAGE 7-9		7-9-91		
GROUND SURFACE ELEVATION:	Г		12" DIA. BORING	DEPTH FROM TOP
GEOLOGIC LOG	DEPTH (FT)		TOP OF GROUND	GROUND (FEET)
FILL, 0-1' gravel 1-9' sandy clay, med to dk gry, rubber like debris @ 5-7'	13333 FIL		BENTONITE-CEMENT GROUT TOP OF BENTONITE	10
SAND, little clay, lt to med gry, in to med, OVM rdg Oppm	9 /// 		POLYMER GROUT  BENTONITE POLYMER GROUT	]
SAND, gry, fine to coarse, loose OVM rdg Oppm	15		6" DIA. PVC, FLUSH JOINT, SCHEDULE 80 RISER PIPE TOP OF BENTONITE -	
		A	PELLET SEAL BENTONITE PELLET SE	EAL .
	S1 :::		- TOP OF FILTER PACK - TOP OF SCREEN	
			6" DIA. TYPE 304 STAINLESS STEEL SCREEN: 0.020" SLOT	
v coarse @ 30' OVM rdg 5ppm @ 32			FILTER PACK SILICA SAND 16-40 SIEVE	
CLAY, lt reddish brn, soft, some gravel	33 :::	<del>   </del>	BENTONITE PELLET SE	<del></del>
NO SCALE (NOTE DEPTH TO	/		BOTTOM OF WELL	— <u>35</u>
DRILLING AND SAMPLING NO	SAMPLIN	<u>.</u>		
OVM READINGS: 0 ppm @		ERVAL (FEET)		
		••		1
	····			

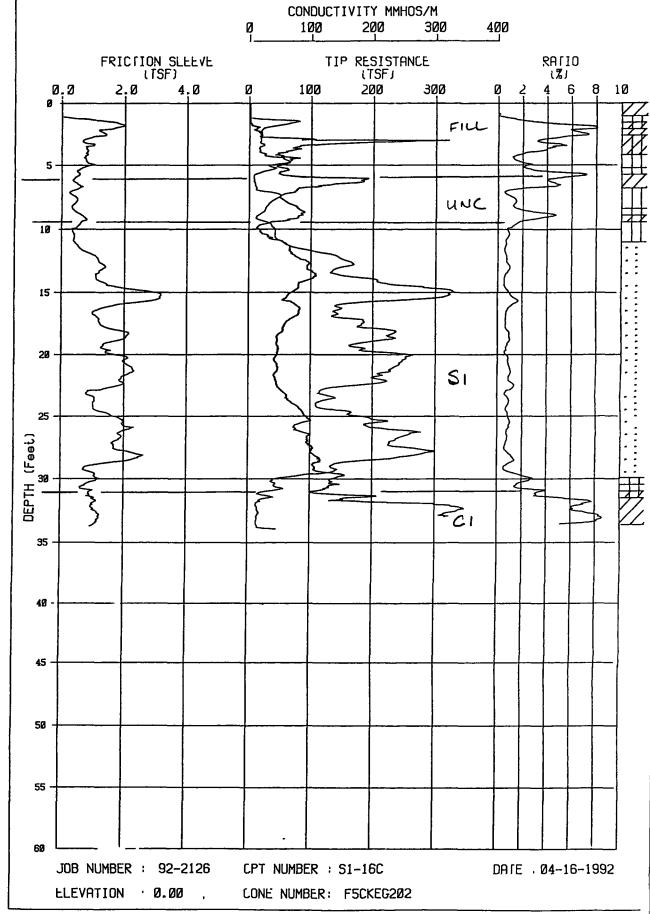




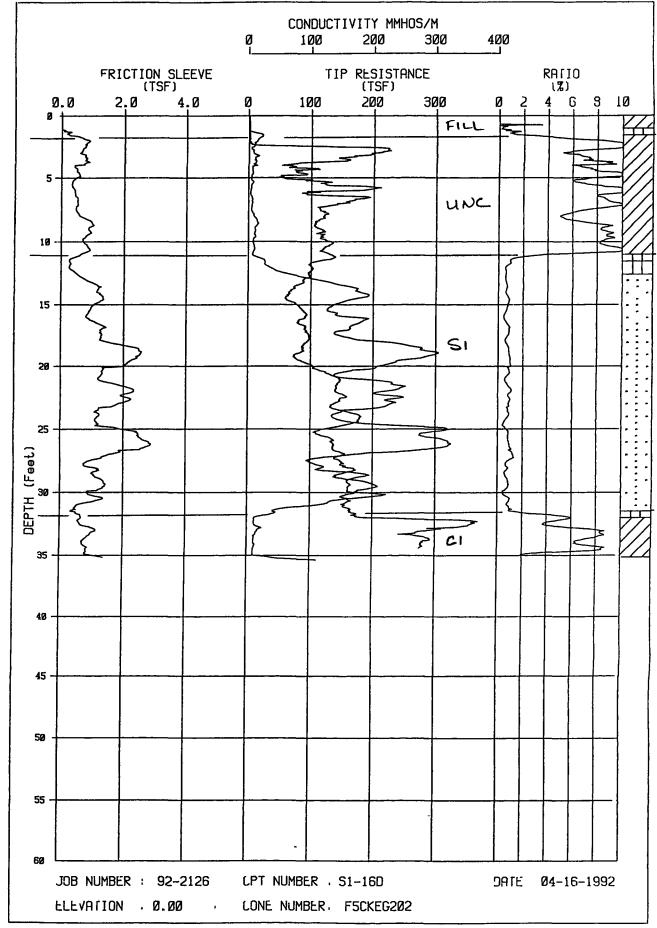
FUGRO GEOSCIENCES, INC



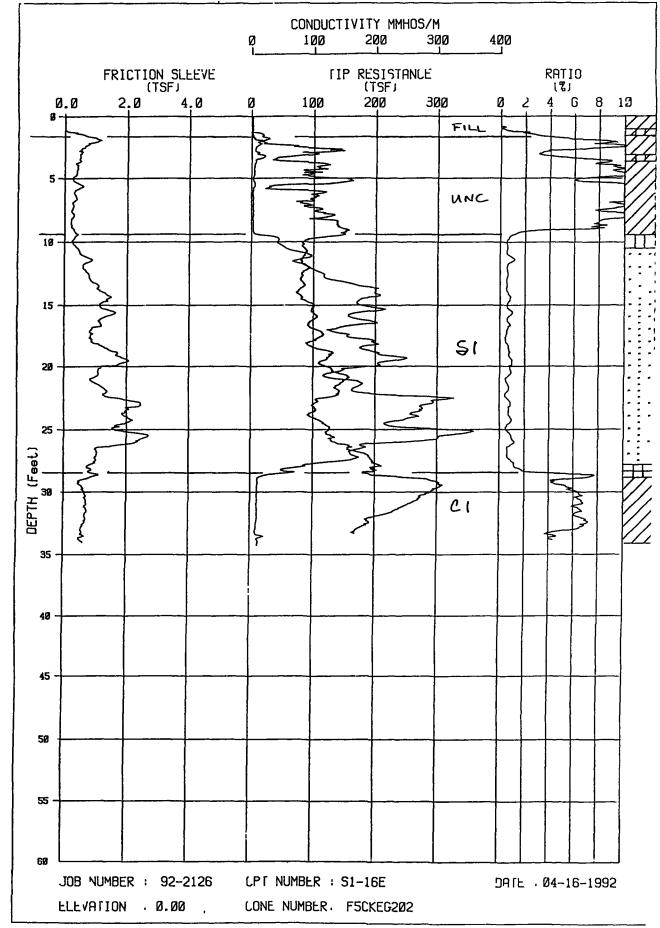
FUCRO GLOSCIENCES, INL



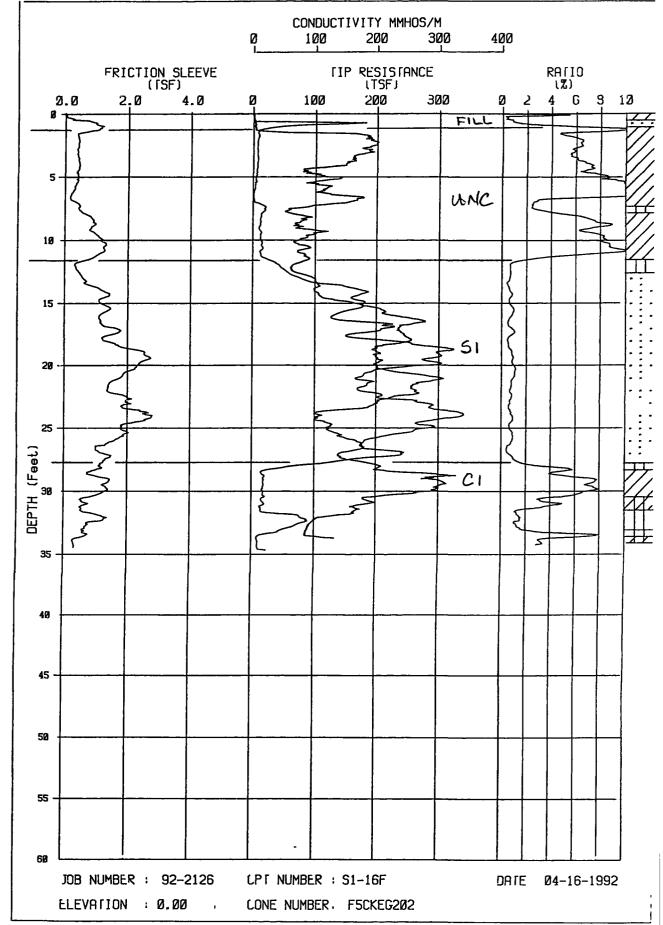
FUCRO GEOSCIENCES INC



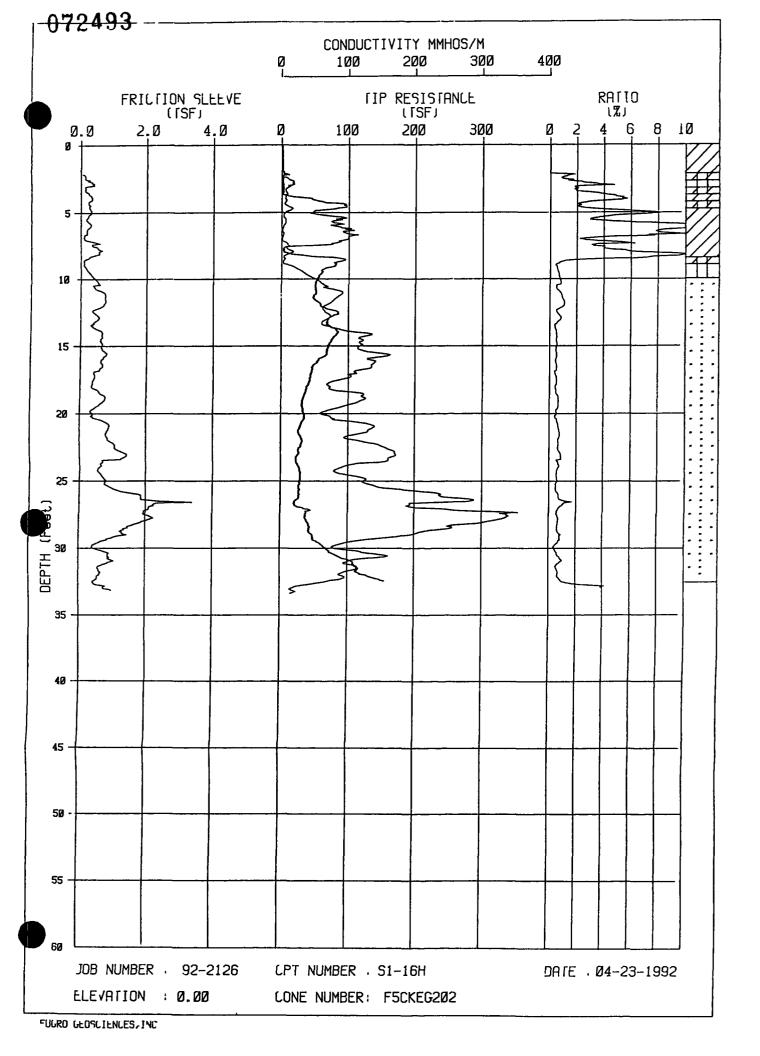
FUGRO GLOSCIENCES, INC

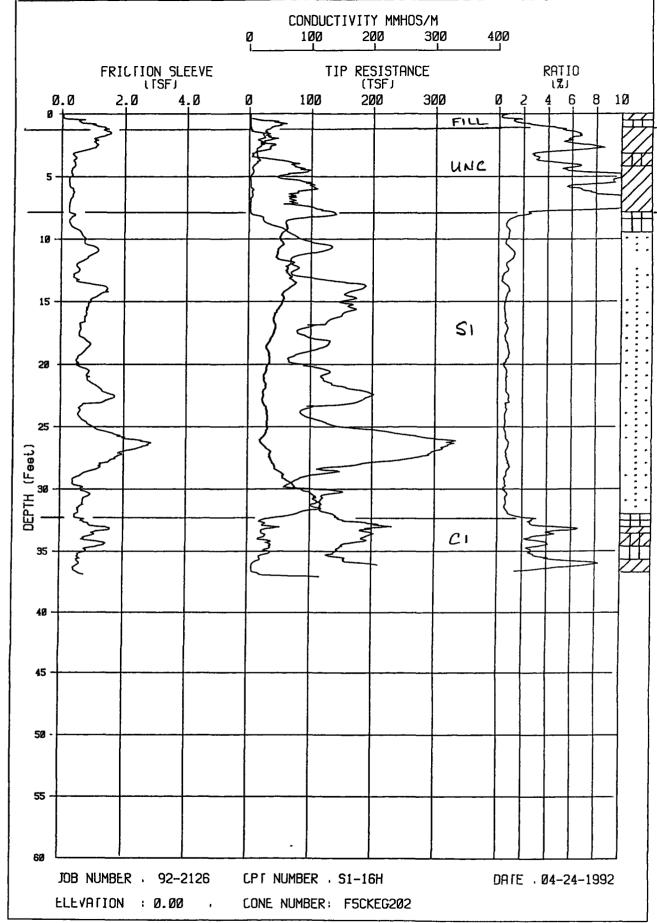


FUGRO GEOSCIENCES, INC

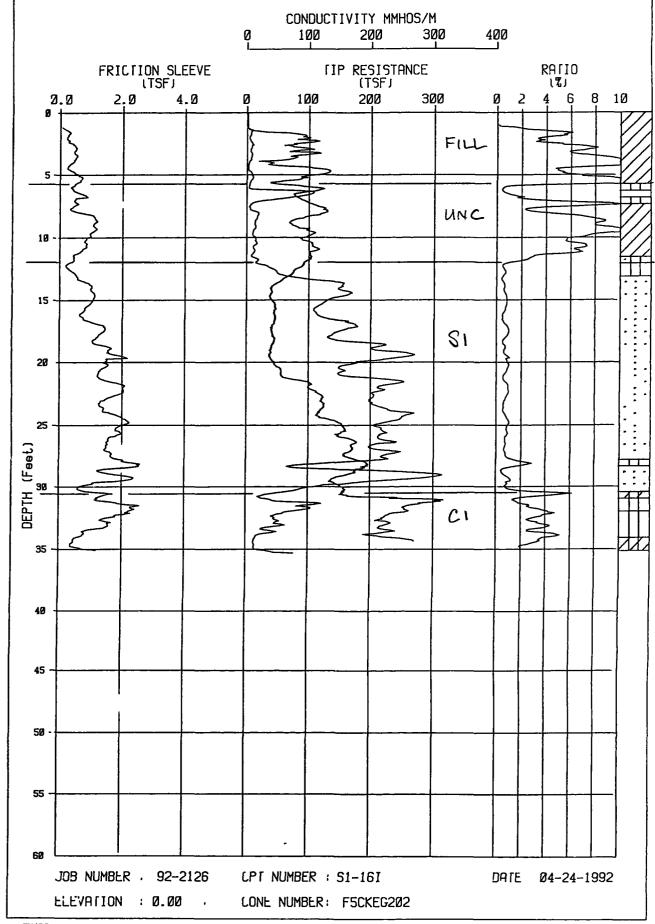


FUGRO GŁOSCIENCES, JNC

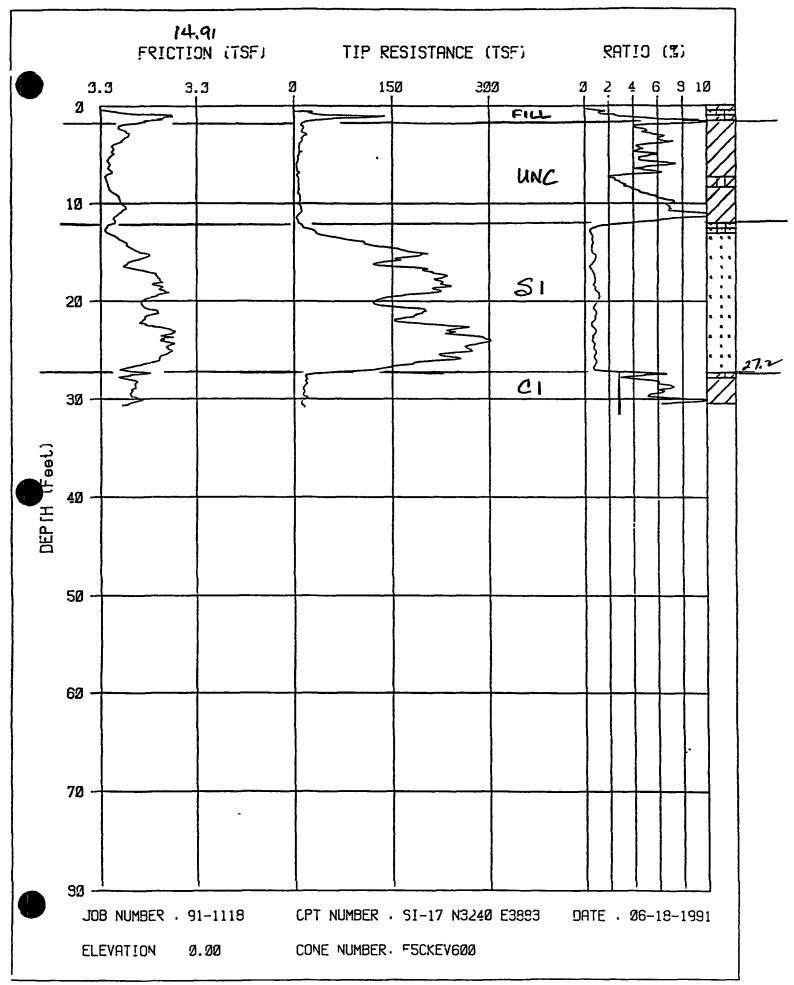




FUGRO GEOSCIENCES, INC

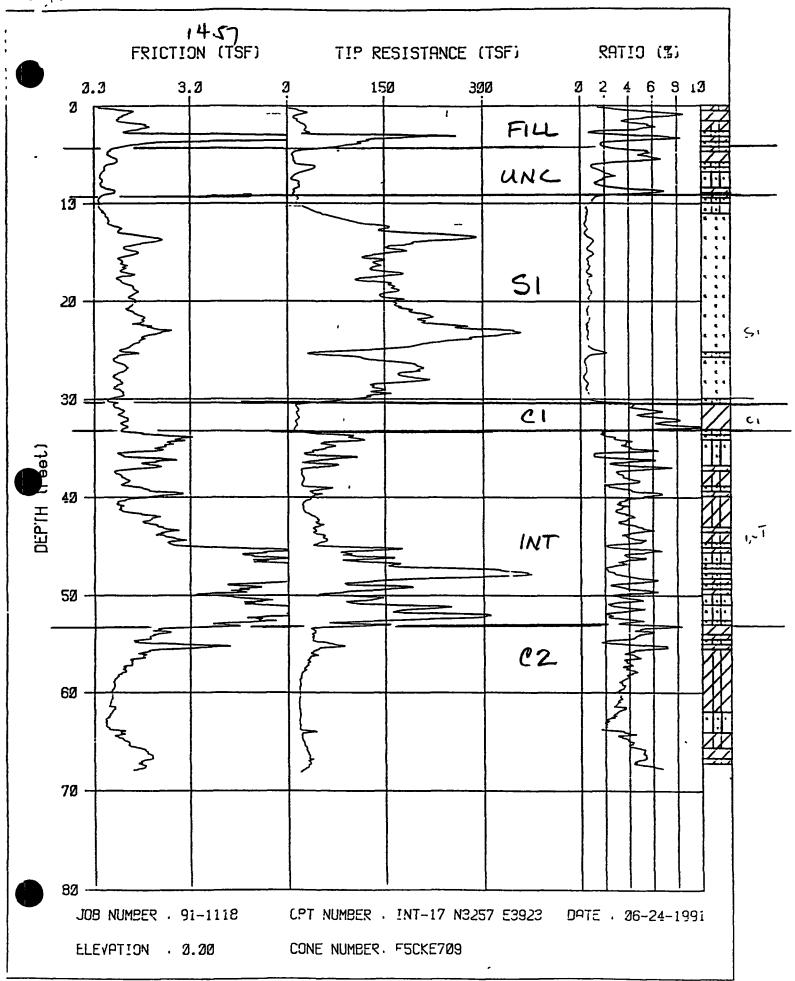


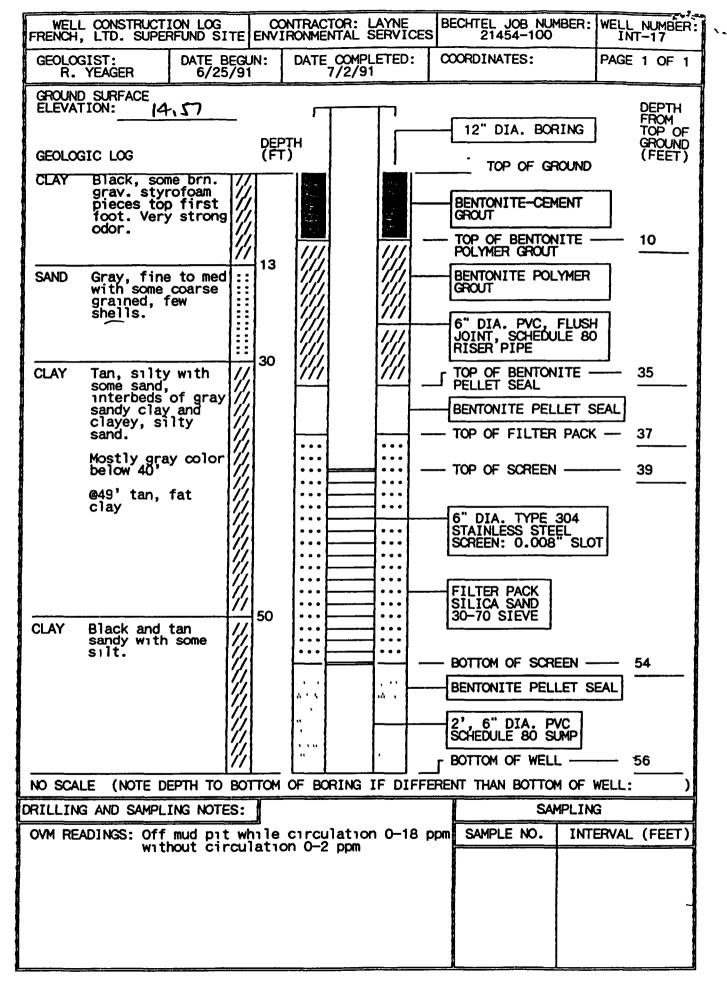
FUGRO GEOSCIENCES, INC



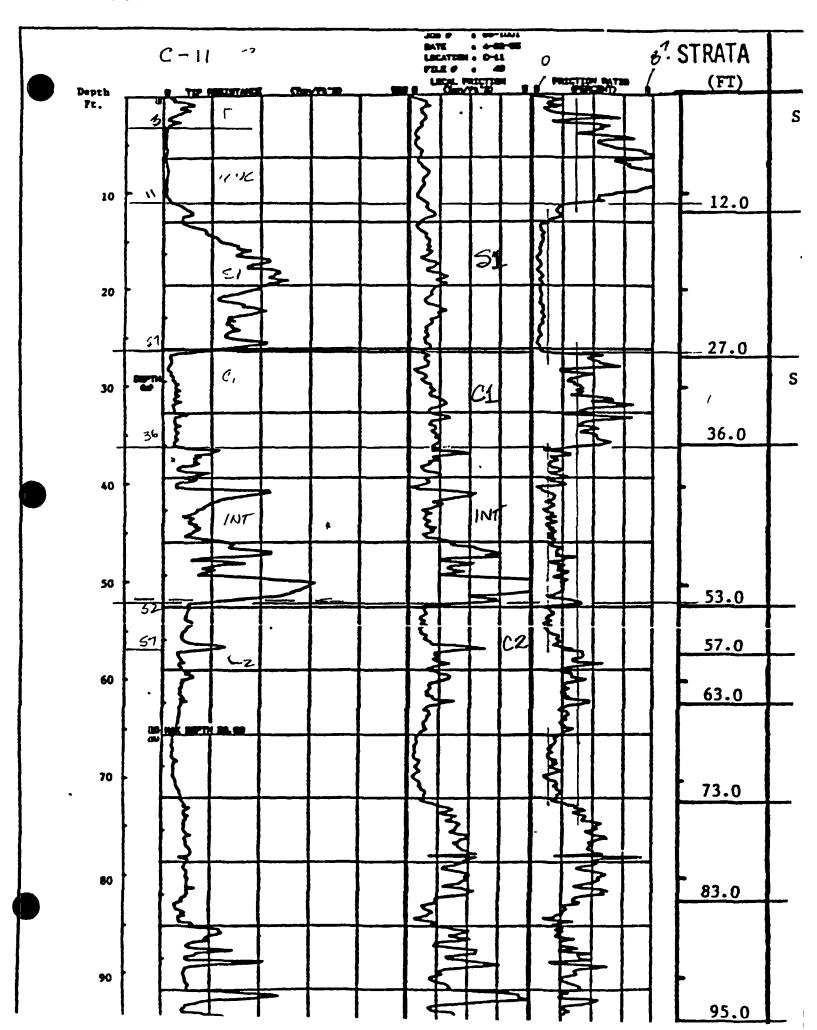
WELL CONSTRUCTION LO FRENCH, LTD. SUPERFUND	G O	ONTRACTOR:	LAYNE SERVICES	BECHTEL JOB NUM 21454-100	MBER: WELL NUMBER: \$1-17
GEOLOGIST: DATE J. T. SMITH 7-	BEGUN: 3-91	DATE COM 7-5		COORDINATES:	PAGE 1 OF 1
GROUND SURFACE ELEVATION: (4.9)  GEOLOGIC LOG  FILL Blk, silty, org. debris  CLAY Charcoal with red blotches, silty, mucky  SAND Tan with black streaks, silty, fine to modium.	DEP (F):			TOP OF GROUT  TOP OF BENTONITE-CEME GROUT  TOP OF BENTONITE POLYMER GROUT  BENTONITE POLY GROUT	GROUND (FEET)  IND  TE 9.0  MER
fine to medium grained, many gas bubbles 26-27' coarse grained with pea gravel				6" DIA. PVC, F JOINT, SCHEDUL RISER PIPE  TOP OF BENTONI PELLET SEAL  BENTONITE PELL  TOP OF FILTER  TOP OF SCREEN  6" DIA. TYPE 3 STAINLESS STEE SCREEN: 0.020"  FILTER PACK SILICA SAND 16-40 SIEVE	TE
CLAY Mottled brown contaminate, blk, slimy	27		n.m.vi	BOTTOM OF SCREE	ET SEAL
SAND Black, clayey cont. froth	<u> </u>		    	2', 6" DIA. PV SCHEDULE 80 SU BOTTOM OF WELL	
NO SCALE (NOTE DEPTH TO DRILLING AND SAMPLING NO		OF BORING	IF DIFFERE	Y	OF WELL: 30.0') MPLING
Highly contaminated wel		required.	-	SAMPLE NO.	INTERVAL (FEET)
	0-9', 5	.0 ppm 9- 85 ppm @ 1		1 2 3	11.5 TO 13.5 19.0 TO 21.0 26.0 TO 28.0









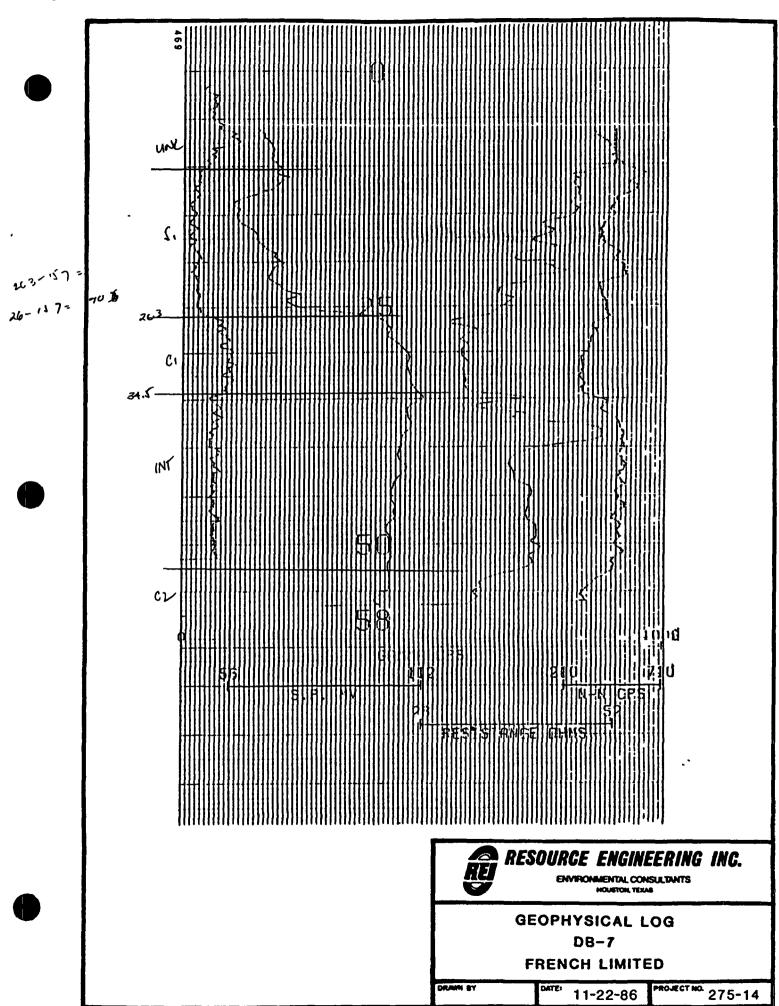




### LITHOLOGIC LOG AND CONSTRUCTION OF DB-7

Client	FRENCH, LTD.	DRILLING AND SAMPLING INFORMATION
Project Name	French, Ltd.	Date Started 7/18/86 Date Completed 7/19/86
Project Location.	Crosby, Texas	Method Rotary Wash Total Depth 58.0
Job No	275-14 Boring NoDB-7	
Logged By	P Mann, Hydrogeologist	
Approved By		
Drilled By	Gulf Coast Coring - Rosenburg, Texas	

DEPTH IN FEET	DESCRIPTION SURFACE ELEVATION 15,7 & F.	STRATUM ELEVATION IN FEET	SAMPLE NO.	SAMPLE TYPE	RECOVERY	GRAPHIC LOG	HNU (IN ppm)	OVA (IN ppm)	SAMPLED FOR CHEM ANALYSIS
l°=	SILTY SAND (FILL?), gray and black; odor detected	2,0	0-2	ST	1.0'	Г	50	Г	
=	SILTY CLAY, gray, massive, moderately firm		2-4	SS	1.0'		-1		'
5-			4-6	ST	1.8'		-1		Ιi
		UNC	6-8	ST	1.9'	ĺ	1.0		
=			8-10	SS	0.3	İ	<1		
10-	SILTY SAND, ranging from fine to coarse grained, gray, poorly sorted,	10.4	10-12	SS	1.4'		-1		1 1
3	g: aded		12-14	SS	1.7'			1.0	
15-3	saturated below 12.0°		14-16	\$S	0.7'		1	6	-
=			16-18	ss	1.1'		<b>~1</b>	80	İ
	18.0° to 22.0° very dark gray, discolored silty sand	SI	18-20	ss	1.3'			100	
20			20-22	ss	1.1'		2	200	1 1
			22-24	SS	1.0'			10	
25	24.5° with pebble sized gravel	26.3	24-26	SS	1.0'		3		
	SILTY CLAY, reddish brown and gray, mottled, massive, very firm	-	26-28	ST	1.5'		3		曰
] =	26.3- ss (	2 3	28-30	ST	1.9'		41		
30 =	·		30-32	ST	1.8'		<1		1
] ]		CI	32-34	ST	1.7'		<b>41</b>		
35-	below 35.6' increasingly sendy, fine grained with grannish gray color		34-36	ST	1.6'		41		
=		38.0	36-38	ST	1.8'		<1		
] =	CLAYEY SAND WITH SILT, very fine grained, greenish gray, moist	4	38-40	ST	1.5'		<1		$\dashv$
40-3		41.0	40-42	SS	1.2*		-1	2	
=	SILTY TO SANDY CLAY, gray, brown and reddish brown, mottled, massive, very firm	1,,,,,	42-44	ST	1.35°		-1	2	$\exists$
45-3		INT	44-46	ST	1.21		41		
	CLAYEY SAND, very fine to fine grained, gray and brown, mottled, dense, moist	1	46-48	ST	1.7'			-1	$\exists$
50-			48-5Q	ST	1.6*			<b>41</b>	
"=	below 52.0' reddish brown color observed		50-52	ST	1.5'			<1	1
=	CLAY TO SILTY CLAY, reddish brown and gray, mottled, massive, very firm	<b>V</b> <sub>53 0</sub>	52-54	ST	1.65			41	
55-		C2	54-56	ST	1.5'			-1	
		58 0	56-58	ST	1 65			-1	$\dashv$
E 60 L	BORING COMPLETED AT 58 O FEET, BOREHOLE WAS GEOPHYSICALLY LOGGED, THEN SEALED WITH CEMENT BENTONITE GROUT.								
	SAMPLER TYPE BOI	ING METHOD							



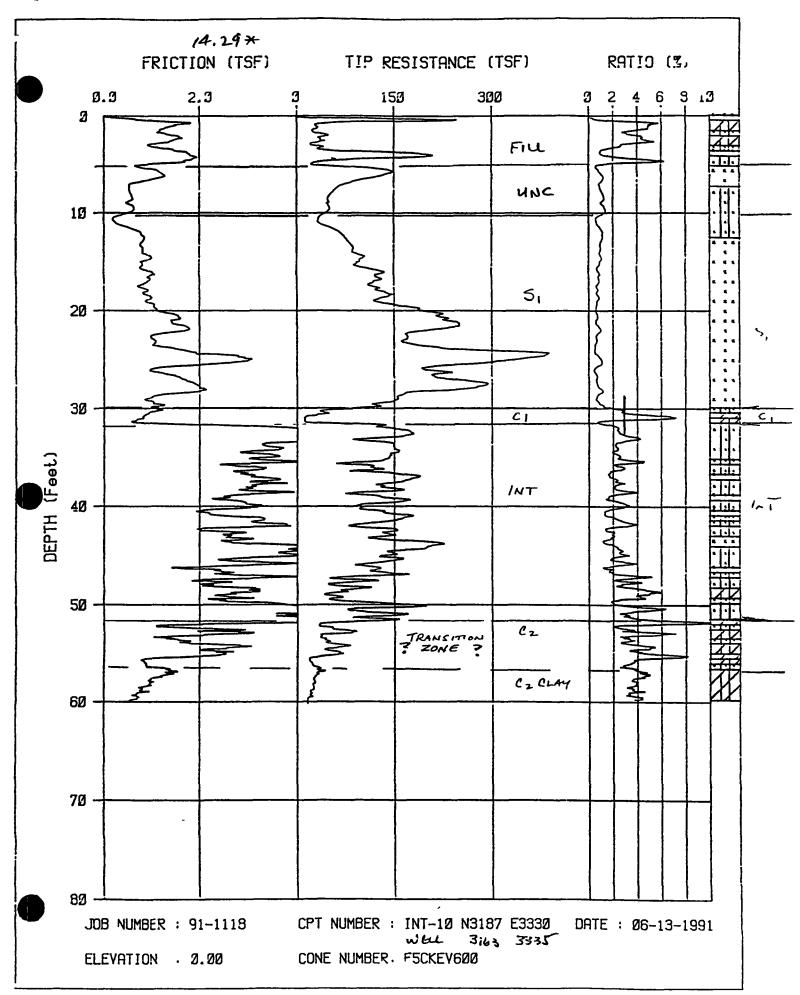
# 

٠,

### Appendix 4

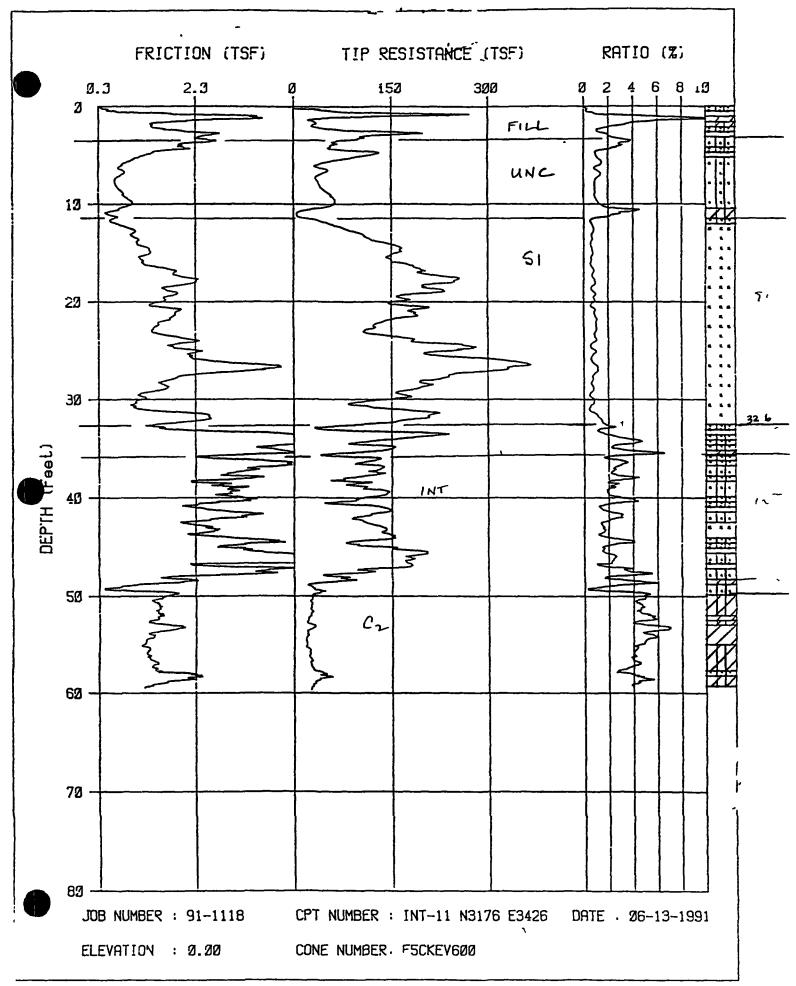
Cone Penetrometer Logs INT-11 Area

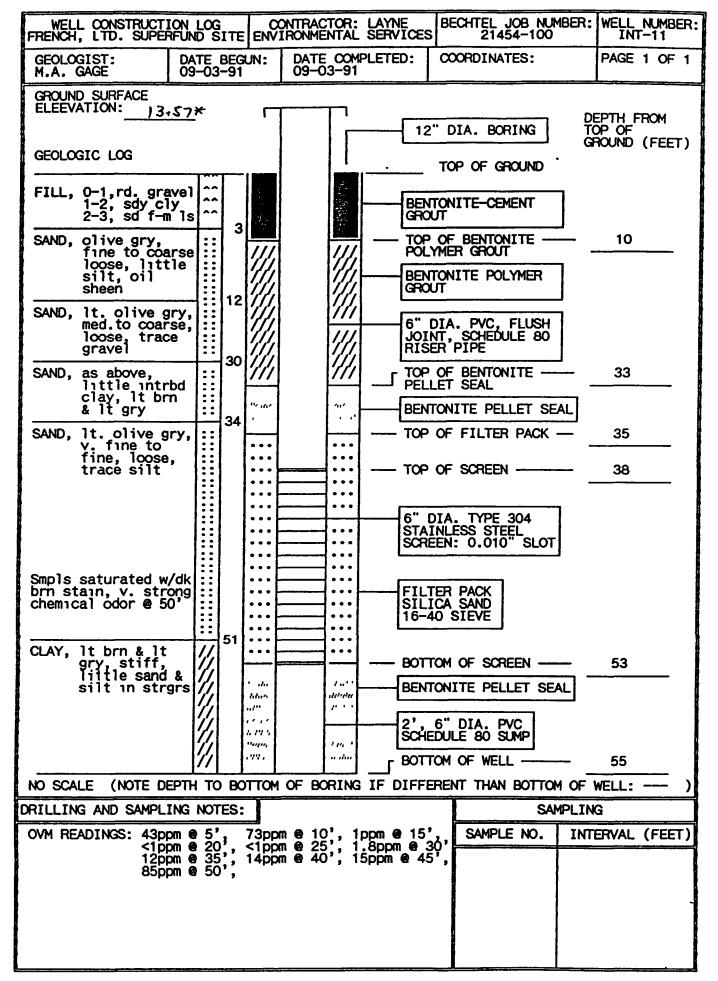
01502436



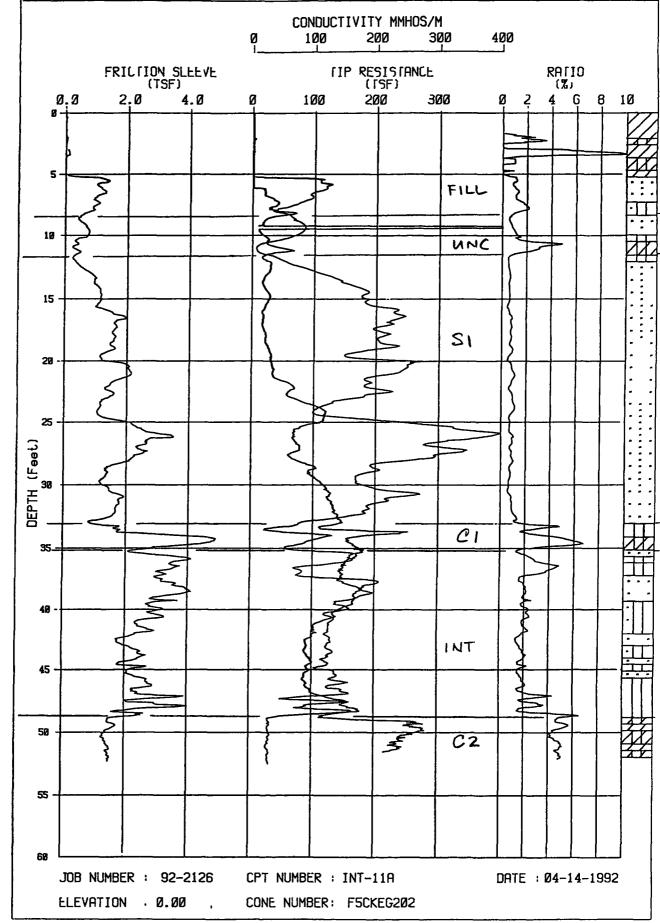
WELL CONSTRUCT! FRENCH, LTD. SUPER	ON LOG FUND SITE EN	CONTRACTOR: /IRONMENTAL	LAYNE SERVICES	BECHTEL JOB NUM 21454-100	IBER: WELL NUMBER
GEOLOGIST: J.T.SMITH	DATE BEGUN: 6-27-91	DATE COMP 7-2-9	PLETED:	COORDINATES:	PAGE 1 OF 1
GEOLOGIC LOG  FILL Road base gailt, org.	rav.	ЕРТН -Т)		12" DIA. BOR	GROUND (FEET)
sand roots.  SAND Tan-gray, m grained, qt loose. 8-13', graye 18-27', bl. staining 27-29 tan, w rnd. pea gr	ed. z, rnd.			BENTONITE-CEM GROUT  TOP OF BENTON POLYMER GROUT  BENTONITE POLY GROUT  6" DIA. PVC, I JOINT, SCHEDU	ITE10
PEA GRAVEL Clayey charcoal with r stains Hly vari well rnd. dk/l petrified wd.	ed iron oo			TOP OF BENTON: PELLET SEAL BENTONITE PELL	33
SAND Grn-gray to tan, fine t grained. pea gravel 49	gray ::			<b>L</b>	PACK — 35.2
SAND & PEA GRAVEL with tan cl iron staine siltstone.	ay and 50			6" DIA. TYPE 3 STAINLESS STEE SCREEN: 0.008	EL I
SAND clayey, sta  SAND Gray green fine to med grained, so staining	to gray :: 51			FILTER PACK SILICA SAND 30-70 SIEVE  BOTTOM OF SCRE	
CLAY Red-brown, gumbo, plas	tic ///	.6)		2', 6" DIA. PV SCHEDULE 80 SL BOTTOM OF WELL	
NO SCALE (NOTE DI DRILLING AND SAMPL		M OF BORING	IF DIFFE		1 OF WELL: 72'
Hole plugged back well.  OVM READINGS: 0.2	to 59' with	1.0 ppm @	50-51'	ng SAMPLE NO.  1 2 3	INTERVAL (FEET 29-30' 30-32' 40-42'
Contaminate found oxide stains and o	in samples @ odor in sampl	53' & 57'. es and cutt	Strong ings.	4	49-50', 51-53



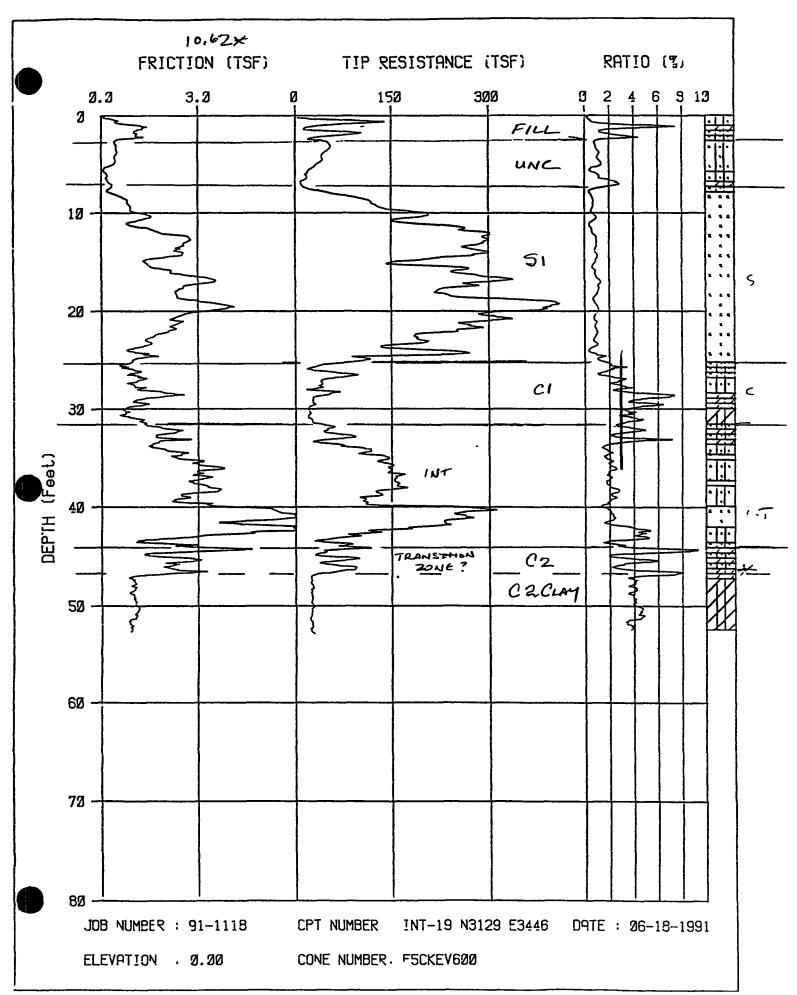


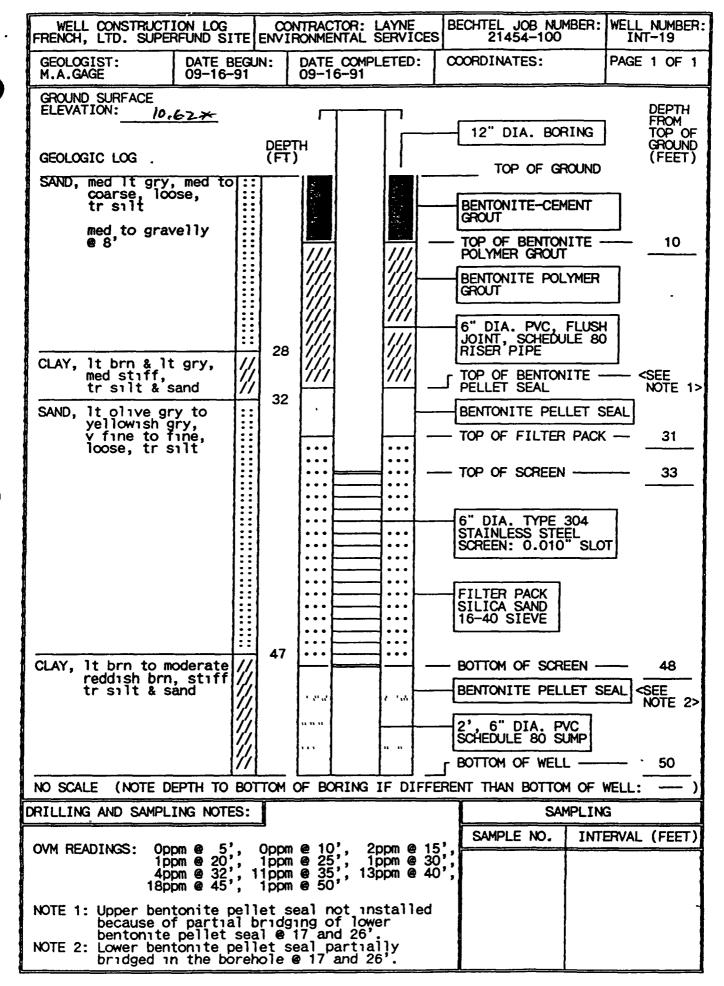




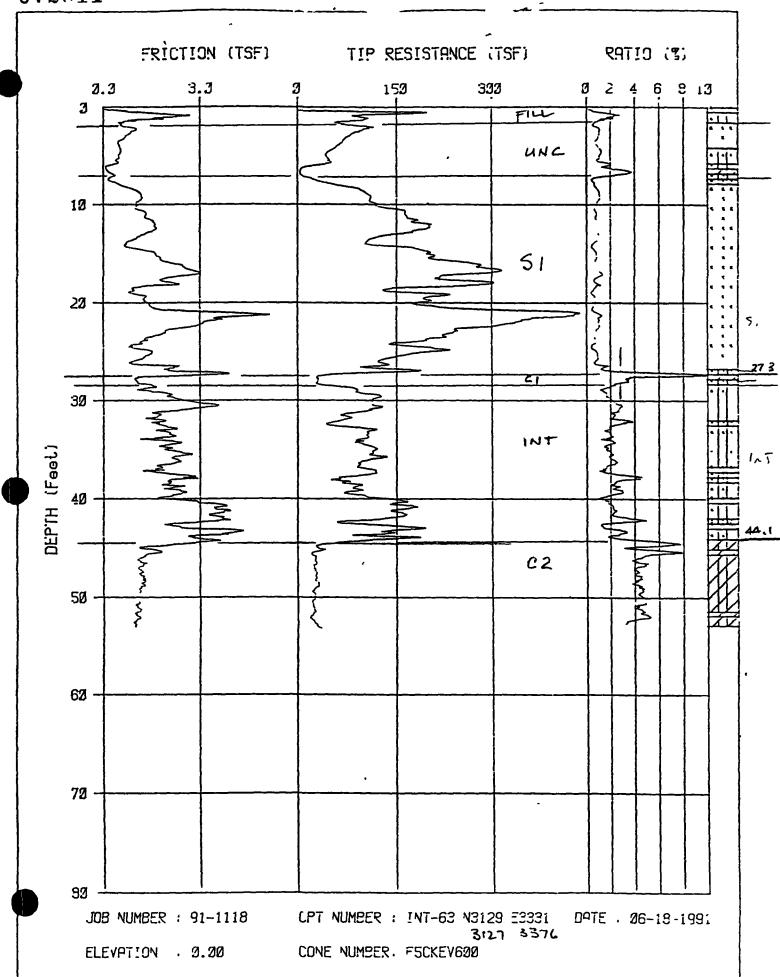


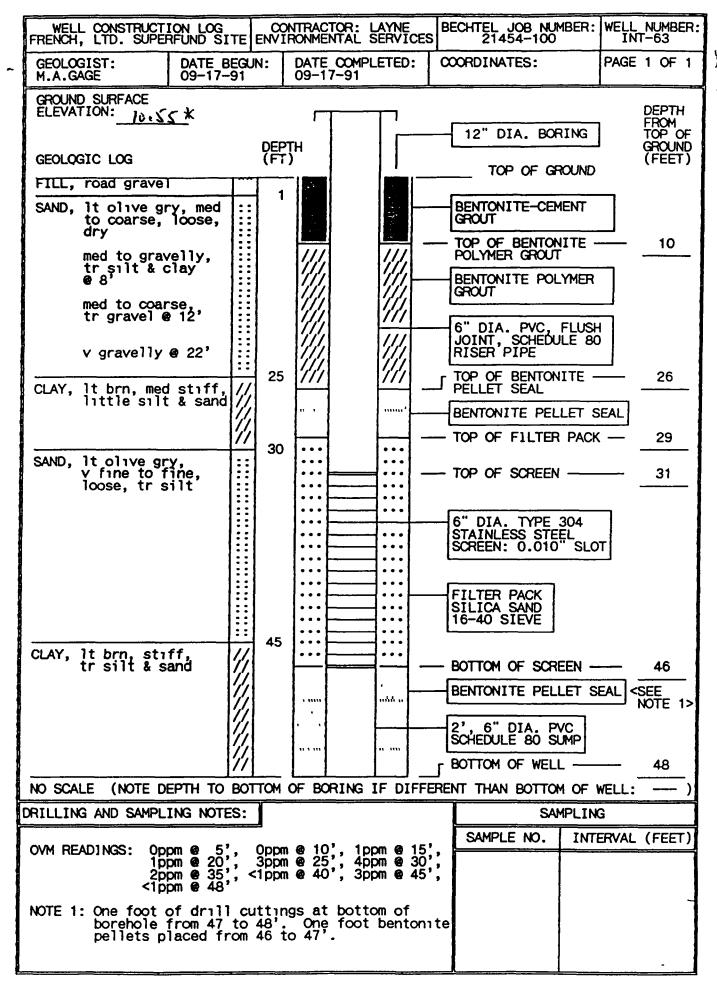
FUCRO GLOSCIENCES, INC



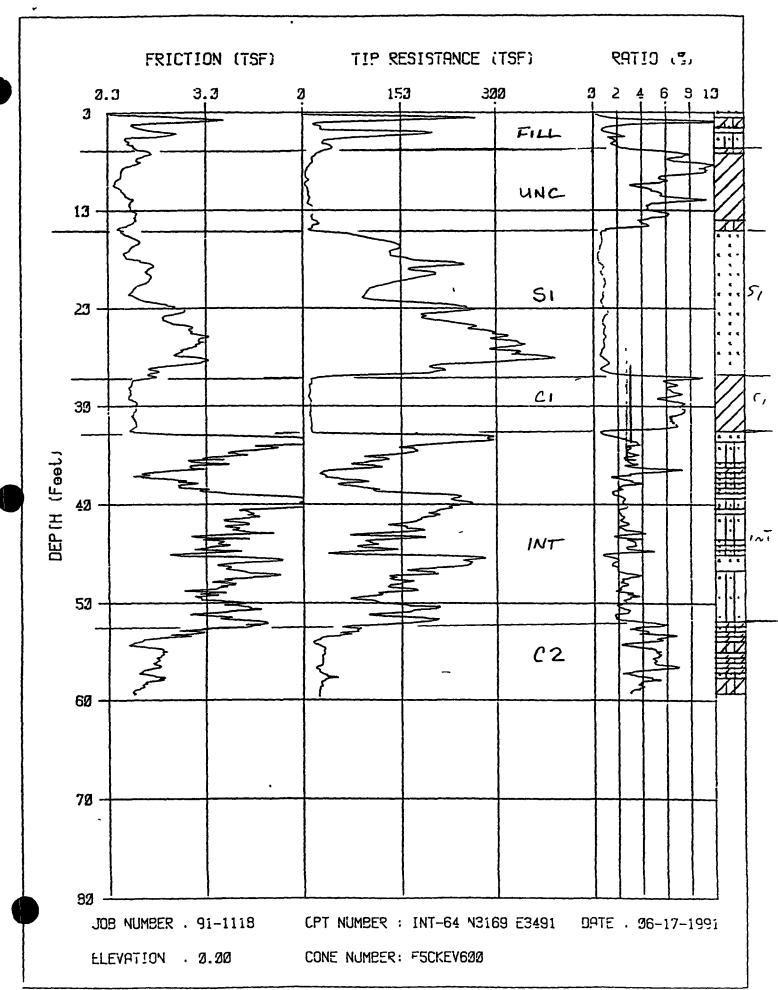










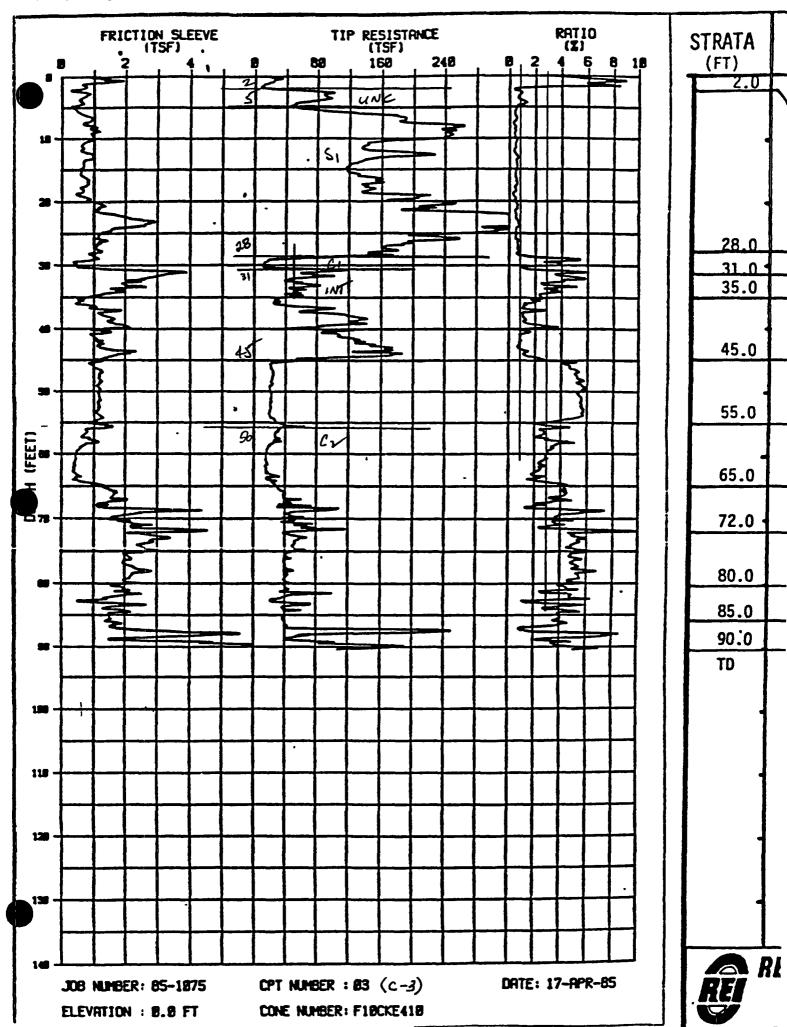


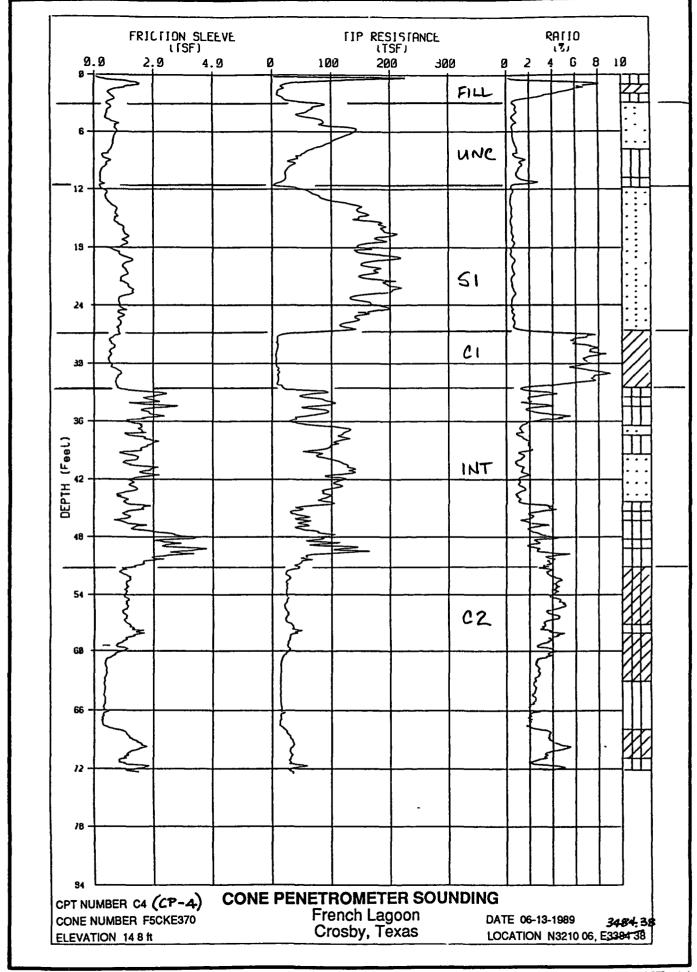
FRENCH, LTD. SUPERFUND SITE ENVIRONMENTAL SERVICES 21454-100 INT-64  GEOLOGIST: DATE BEGUN: DATE COMPLETED: COORDINATES: PAGE 1 OF 1  GROUND SURFACE ELEVATION: 13.87*  GEOLOGIC LOG  FILL, roadbase  SAND, sity, clyey, brn w/bik stn & oily size.  SAND, cs-v cs, blk stn, loose washes, blk blobs of cont 15-17 cs, frly cln, tan, clrlt grns, rnded, med-17-24' as above, cs-v cs @ 22' tr pea grv! blk cont. pea grv! blk cont. pea grv! blk cont. pea grv! blk cont. pea gry! slk cont. pea gry! bls blk cont. pea gry! slk cont. pea gry! slk cont. pea gry! slx slk cont. pea gry! slx slk cont. pea gry! slx slk cont. pea gry! slx slk cont. pea gry! slx slk cont. pea gry! slx slk cont. pea gry! slx slk cont. pea gry! slx slk cont. pea gry! slx slk cont. pea gry! slx slk cont. pea gry! slx slx slk cont. pea gry! slx slx slk cont. pea gry! slx slx slx slx slx slx slx slx slx slx										
ELEVATION: 13.67*  GEOLOGIC LOG  GEOLOGIC LOG  GEOLOGIC LOG  GEOLOGIC LOG  FILL, roadbase  SAND, slty, clyey, brn w/blk stn & oily scum w/blk stn & oily scum sense, blk blobs of cont 15-17 cs, frilt grn. ender, meding from 17-24 as above, cs-v cs & 22', rr 21-24', rd-blk blotches 24-28' clyey blbs blk cont, pea grvl blk cont gry stn, loose 25-28!  CLAY, tan-rddsh, tight, dnse, gry strks slckensd brks, plastic, dk-rd 30' // SILT, tight, dnse, it gry, w/ocher stns, sickensd brks, plastic, dk-rd 30' // SILT, fr, wet, wife od; v aromtc, it berge  SILT, fn, crmy bge, frly mill sem 1-31 blick, pliable 50-52' crmy bge, frly sem 1-31 blick, pliable 50-52' crmy bge, mill sem 1-31 blick pliable 50-52' crmy bge, mill sem 1-31 blick pliable 50-52' crmy bge, mill sem 1-31 blick pliable 50-52' crmy bge, mill sem 1-31 blick pliable 50-52' crmy bge, mill sem 1-31 blick pliable 50-52' crmy bge, mill sem 1-31 blick pliable 50-52' crmy bge, mill sem 1-31 blick pliable 50-52' crmy bge, mill sem 1-31 blick pliable 50-52' crmy bge, mill sem 1-31 blick pliable 50-52' crmy bge, mill sem 1-31 blick pliable 50-52' crmy bge, mill sem 1										
SAND, slty, clyey, brn w/blk stn & oily scum  SAND, cs-v cs, blk stn, loose, washes, blk blobs of cont 15-17 cs, frly cln, tan, cir-lt grns, rnded, med- fn e 16-17' 17-24 as above, cs-v cs e 22', tr pea grv1, blk cont 21-24', rd-blk blotches 24-28' clyey blbs blk cont, pea grv1 gry stn, loose e 25-28'  CLAY, tan-rddsh, tight, dnse, gry strks slckensd brks, plastic, dk-rd 30'  SILT, tight, dnse, lt gry, w/ocher stns, riable 33.5-41' 10% cly  SAND, slty, fn, wet, wife od; v aromtc, it-berge SILT, fn, crmy bge, frly soft, wet 43  SILT, fn, crmy bge, frly soft, wet semi-dnse, rd-org ocher blotches  2', 6" DIA, PVC  BENTONITE CEMENT FOO OF BENTONITE 9  Colymen GROUT  TOP OF BENTONITE POLYMER GROUT  FOLYMER GROUT  TOP OF BENTONITE POLYMER BENTONITE POLYMER GROUT  TOP OF BENTONITE  9  Colymen GROUT  TOP OF BENTONITE  13  BENTONITE CEMENT  TOP OF SCREEN  TOP OF FILTER PACK STAINLESS STEEL SCREEN: 0.010" SLOT  41  43  SILT, fn, crmy bge, frly soft, wet 43  SILT, plable 50.52 crmy bge, wine od, frla, semi-dnse, rd-org ocher blotches  TOP OF SCREEN  T										
SAND, slty, fn, wet, semi-dase, rd-org ocher blotches   SILT, fn, crmy bge, frly soft, wet 48-50', 50% cly, rddsh gry, soft, wet 48-50', 50% cly, rd-org ocher blotches   Soft, wet wine od; rd-org ocher blotches   Soft, semi-dase, rd-org ocher blotches   Semi-dase, rd-org ocher blotches   Silica, semi-dase, rd-org ocher blotches   Silica, semi-dase, rd-org ocher blotches   Silica, semi-dase, rd-org ocher blotches   Silica, semi-dase, rd-org ocher blotches   Silica, semi-dase, rd-org ocher blotches   Silica, prime of semi-dase, rd-org ocher blotches   Silica, prime of semi-dase, rd-org ocher blotches   Silica, prime of semi-dase, rd-org ocher blotches   Silica, prime of semi-dase, rd-org ocher blotches   Silica, prime of semi-dase, rd-org ocher blotches   Silica, prime of semi-dase, rd-org ocher blotches   Silica, prime of semi-dase, rd-org ocher blotches   Silica, prime ocher blotches										
Diotoches 24-28' clyey blbs blk cont, pea grv1 gry stn, loose 6 25-28'  CLAY, tan-rddsh, tight, dnse, gry strks slckensd brks, plastic, dk-rd 30'  SILT, tight, dnse, lt gry, w/ocher stns, friable 33.5-41' 10% cly  SAND, slty, fn, wet, it-berge  SILT, fn, crmy bge, frly soft, wet 41  SILT, fn, crmy bge, frly soft, wet 43  SILT, fn, crmy bge, frly soft, wet 45-05' 50% cly, rddsh gry, soft, slick, pliable 50-52 crmy bge, wine od, fria, semi-dnse, rd-org ocher blotches  SILT, for the following bge, wine od, fria, semi-dnse, rd-org ocher blotches  SILT, for the following bge, wine od, fria, semi-dnse, rd-org ocher blotches  SILT, for the following bge, wine od, fria, semi-dnse, rd-org ocher blotches  SILT, for the following bge, wine od, fria, semi-dnse, rd-org ocher blotches  SILT, for the following bge, wine od, fria, semi-dnse, rd-org ocher blotches  SILT, for the following bge, wine od, fria, semi-dnse, rd-org ocher blotches  SILT, for the following bge, wine od, fria, semi-dnse, rd-org ocher blotches										
CLAY, tan-rddsh, tight, dnse, gry strks slckensd brks, plastic, dk-rd 30'  SILT, tight, dnse, lt gry, w/ocher stns, friable 33.5-41' 10% cly  SAND, slty, fn, wet, wine od; v aromtc, lt berge  SILT, fn, crmy bge, frly soft, wet 48-50' 50% cly, rddsh gry, soft, wet 50-52' crmy bge, wine od; fria, sem1-dnse, rd-org ocher blotches  SENTON OF SCREEN 37  TOP OF SCREEN 37  6" DIA. TYPE 304  STAINLESS STEEL SCREEN: 0.010" SLOT  FILTER PACK SILICA SAND 16-40 SIEVE  BENTONITE PELLET SEAL  2'. 6" DIA. PVC										
SILT, tight, dnse, it gry, w/ocher stns, iiiiiiiiiiiiiiiiiiiiiiiiiiiiiiiiiiii										
SAND, slty, fn, wet, wine od; v aromtc, it beige  SILT, fn, crmy bge, frly soft, wet 48-50' 50% cly, rddsh gry, soft, slick, pliable 50-52; crmy bge, wine od; fria, semi-dnse, rd-org ocher blotches  FILTER PACK SILICA SAND 16-40 SIEVE  BENTONITE PELLET SEAL										
48-50' 50% Cly, rddsh gry, soft, slick, pliable 50-52 crmy bge, wine od, fria, semi-dnse, rd-org ocher blotches 50' 50% Cly, rd-org ocher blotches 52', 6" DIA, PVC										
CLAY, solid, dnse, tough // 52 SCHEDULE 80 SUMP BOTTOM OF WELL 54										
NO SCALE (NOTE DEPTH TO BOTTOM OF BORING IF DIFFERENT THAN BOTTOM OF WELL: )  DRILLING AND SAMPLING NOTES:  SAMPLING										
During drilling a considerable amount of blk blobs of contaminate (gooey-thick) was trapped within the sands. In air/sunlight, most blobs oxidized completely.  OVM READINGS: 3ppm @ 15-17', 3ppm @ 24-26', 50ppm @ 27-29', 15ppm @ 31-33', 48ppm @ 41-43', 32ppm @ 50-52'  SAMPLE NO. INTERVAL (19 24-26', 24-26', 31-33', 31-33', 32ppm @ 24-26', 31-33', 32ppm @ 50-52'										



MONITORING WEI	L LOG ORFUND SITE ENV	ONTRACTOR:	LAYNE SERVICE	BECHTEL JOB NUMBER 21454-100	: WELL NUMBER:
GEOLOGIST: M.A.GAGE	DATE BEGUN: 09-05-91	DATE COM 09-09-91	PLETED:	COORDINATES:	PAGE 1 OF 1
GEOLOGIC LOG FILL, Oyster she	(F			12" DIA. BORING	- GROUND . (FEET)
SAND, it olive gr brn, med, i SAND, as above, t & clay	∞se			BENTONITE-CEMENT GROUT  TOP OF BENTONITE POLYMER GROUT  BENTONITE POLYMER GROUT  6" DIA. PVC, FLUS	3
CLAY, med lt gry, stiff, some SAND, lt olive gr	/// 29	1 31111		JOINT, SCHEDULE ERISER PIPE  TOP OF BENTONITE PELLET  BENTONITE PELLET  TOP OF FILTER PACE	
y fine to f loose, tr s	ine,			TOP OF SCREEN —  6" DIA. TYPE 304 STAINLESS STEEL SCREEN: 0.010" SL  FILTER PACK SILICA SAND 16-40 SIEVE	31
CLAY, 1t brn, sti	11 12			BENTONITE PELLET  2', 6" DIA. PVC SCHEDULE 80 SUMP  BOTTOM OF WELL —	SEAL 48
DRILLING AND SAMPLE	Opm @ 5', <1p	of BORING  pm @ 10', pm @ 25', pm @ 40',	6ppm @	15' 30'	
Slight chemical to	o mud pit @ 27	to 48'.			







### ENSR CONSULTING & ENGINEERING

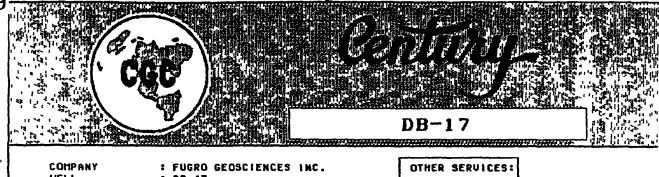
# SUBSURFACE EXPLORATION

Client . French Limited Task Group
Project Name : French Limited Design Phase
Project Location : Crosby, Texas
Job Number : 2870—017 Boring No : DB-17
Logged By : Cameron Teller
Approved By : Hoyt Clark
Prilled By : Fugra Geosciences, Inc. DRILLING AND SAMPLING INFORMATION
Date Started: 9-18-90 Date Completed: 9-19-90
Method: Mud Rotary Total Depth: 54 ft
WELL COMPLETION INFORMATION
Screen Dia: Length:
Siot Size:
Casing Dia: Length:

LITHOLOGIC LOG OF DB-17

r	 				<u> </u>													<del></del>								,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,
	 4.1.1		ي است ا	ш	ئ ىلى		, ,,,,,	<b>Ֆ</b> Լու	الللا	<u>چي</u> سالت	ш	, ,,,,,,,	5 1		25 	; 	8 	; 	5 1		5 		ى دىلى			depth In feet
SAMPLER TYPE	C2	ed-brown clay Blacky texture, slickensided, with sit partings	0		Interbedded stiff greenish—gray silty clay and loose gray fine sand [Moderate to strong odor.]		Z			No CIT		Dense gray clavey coarse sand, some fine grayel.	ome medium and coarse sand 28—29 ft		-little gravel 23-28 ft		C					Loose brown medium sand, with roots 6-10 ft.	6.5		Fill; rock fragments, sand-änd clay.	DESCRIPTION SURFACE ELEVATION :/ 141,5
	 20	19	18	17	16	5	14	<u>ت</u>	12	11	10	9	8		7		6		Ch		4	U	2	-		SAMPLE NO
_	 ST	ST	ST	ST	ST	ST	ST	ST	ST	ST	ST	ST	ST		ST		TS		ST	ļ	ST	ST	श	ST		SAMPLE TYPE
BORN	 				<u> </u>	ļ		<u> </u>							ļ											
E E	 				<u> </u>	<u> </u>		<u> </u>		_					ļ											CUCVIICAL
BORING METHOD	 1	<u> </u>	<u> </u>	>	<u>\</u>	\	5	<u> </u>	<u> </u>	<u>\</u>	-	<u> </u>	\		1		<u>\</u>		<u>\</u>		\	<u> </u>	<u>\</u>	<u> </u>	_	CHEMICAL ANALYSIS GEOTECH ANALYSIS
	 1		1		4				<u> </u>			4	L	L	1		_		<u> </u>	L	<u> </u>					ANALYSIS
		III		777,	777						<u>=</u>	<u>N</u> .	•		····	<u>.                                    </u>	<u> </u>			:				Ш	Ш	
	 														G	ROUT	TO S	URFAC	Έ 							BORING COMPLETION
															1		1									

SECTION



COMPANY : FUGRO GEOSCIENCES INC.

HELL : DB-17

LOCATION/FIELD : FRENCH LTD.

COUNTY : HARRIS

STATE : TEXAS

PERMANENT DATUM : GL DATE : 09/19/90 ELEVATIONS DEPTH DRILLER : 54 J.H. ELEV. PERM. DATUM: GL KB : GL LOG BOTTOM LOG MEASURED FROM: GL 53 70 DF : GL LOG TOP -4.00 DRL MEASURED FROM: GL GL : GL

TOWNSHIP

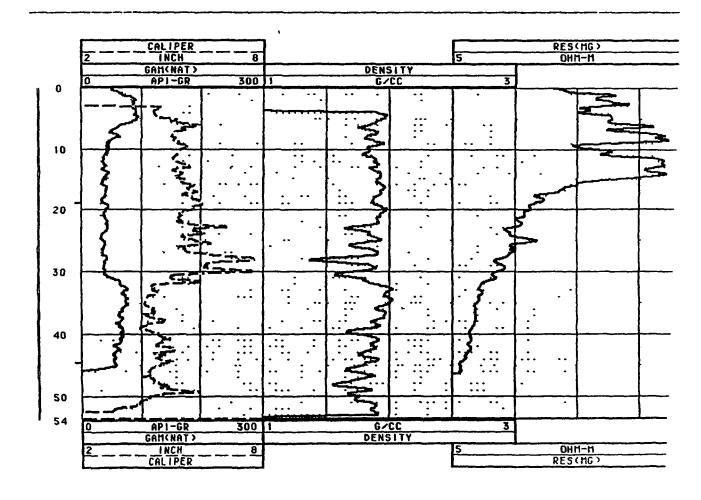
RANGE :

CASING DRILLER: 00 LOGGING UNIT: 9002
CASING TYPE: 00 FIELD OFFICE: TULSA.OK.
CASING THICKNESS: 00 RECORDED BY: I.F.ALANIZ

BIT SIZE : 3.874 BOREHOLE FLUID : H2O/GEL FILE : ORIGINAL MAGNETIC DECL. : 5.8 RM : O TYPE : 9030AA HATRIX DENSITY : 271 RM TEMPERATURE : O LOG : 1 FLUID DENSITY : 1.1 HATRIX DELTA T : 57 PLOT - COLTER :

FLUID DENSITY: 1.1 MATRIX DELTA T: 57 PLOT . COTTER 0
NEUTRON MATRIX: SANDSTONE FLUID DELTA T: 189 THRESH: 50000
REMARKS:

ALL SERVICES PROVIDED SUBJECT TO STANDARD TERMS AND CONDITIONS





### ENSR CONSULTING & ENGINEERING

### SUBSURFACE EXPLORATION

### LITHOLOGIC LOG OF DB-18

DRILLING AND SAMPLING INFORMATION Client: French Limited Task Group

Client: French Limited Task Group
Project Name: French Limited Design Phase
Project Location: Crosby, Texas
Job Number: 2870—017 Boring Na: D
Logged By: Cameron Teller
Approved By Hoyt Clark
Drilled By: Fugro Geosciences, Inc.

Boring Na : DB-18

Date Storted: 9-17-90 Date Completed: 9-18-90
Nethod: Mud Rotary Total Depth. 54 ft
WELL COMPLETION INFORMATION
Screen Dia: Langth:

Slot Size : Type: Casing Dia: Longth :

	by: regro departments, inc. Cosing Dig :			Congu					
DEPTH IN FEET	DESCRIPTION SURFACE ELEVATION: (14,5)	SAMPLE NO	SAMPLE		CHEMICAL	GEOTECH ANALYSS	CRAPHIC LOC	COMPLETION	
	Fill; clay and sand, with roots.						$\blacksquare$		П
5		1 25	ST ST		1				-
1	6.5	3	ST		1		Ш		łł
10	Loose tan to gray fine sand, little to some silt. Slight odor.	4	ST		1				
11111	-black 13-14 ft	5	SS		1				
15		Ĕ	33		-			SURFACE	
20-	5)	6	ST		1	1		2	
25-	-moderate to strong odor below 23 ft	7	ST		1		-	GROUT	
30-1	-below 28 ft loose gray and white medium to coarse sand	8	ST		1	1			
1	-some fine to coarse gravel 30-32 ft -/7,5	9	SS ST		1				
35	Dense green-gray silt, moderate odor.	10 11	ST		1				
11111	•	12 13	ST ST		1	1			
40-1	-black 40.4-40.6 ft, 41 5-42 5 ft, 43.7-43 9 ft	14	ST		1		1111		-
45	INT		ST		1				
45 -	1		ST ST		1				
50-	-fine sandy below 48 ft	18	ST		1				
7 4	$\sqrt{-37.5}$ Very stiff red-brown clay, blocky texture, slickensided. 0.2	19			1		鲗		
55 – 55 –	Very stiff red-brown clay, blocky texture, slickensided. C2	20	ST	-	/	<b>✓</b>			
111111									
ce 14	SAMPLER TYPE  PREN SPIT SPINN BC - BOCK CODE  USA - HOLL	nw <		BORING M	ETHO		0.0	HVING C	



COMPANY : FUGRO GEOSCIENCES INC.

HELL : DB-18

LOCATION/FIELD : FRENCH LTD.

COUNTY : HARRIS

STATE : TEXAS
SECTION :

DATE : 09/18/90 PERMANENT DATUM : GL **ELEVATIONS** DEPTH DRILLER : 54 J.M. : GL ELEV. PERM. DATUM: GL KΘ LOG BOTTOM 53.50 LOG MEASURED FROM: GL DF : GL LOG TOP -3.40 DRL MEASURED FROM: GL GL : GL

TOWNSHIP

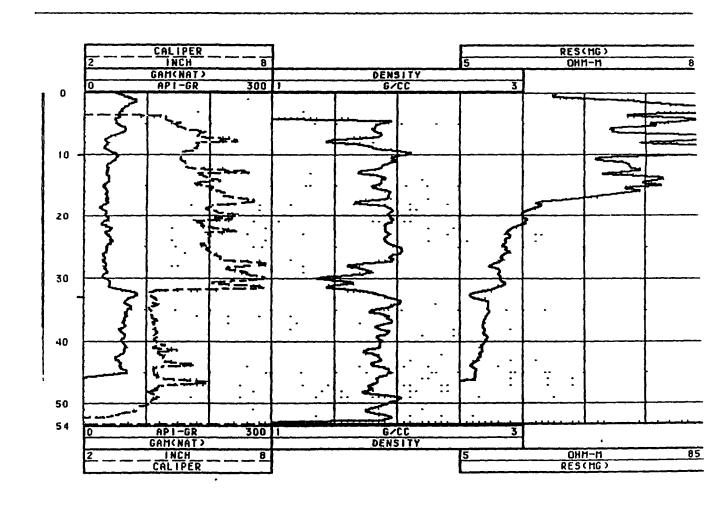
RANGE :

CASING DRILLER: 00 LOGGING UNIT : 9002
CASING TYPE: 00 FIELD OFFICE : TULSA.OK.
CASING THICKNESS: 00 RECORDED BY : I.F.ALANIZ

BIT SIZE : 3.874 BOREHOLE FLUID : H20/GEL FILE : ORIGINAL MAGNETIC DECL. : 5.8 RH : 0 TYPE : 9030AA RM TEMPERATURE MATE IX DENSITY : 271 : 0 LOG : 8 FLUID DENSITY : 1.1 MATRIX DELTA T : 57 PLOT : COTTER 0 FLUID DELTA T NEUTRON MATRIX : SANDSTONE : 189 THRESH: 50000

REMMRKS :

### ALL SERVICES PROVIDED SUBJECT TO STANDARD TERMS AND CONDITIONS



### **ENSR CONSULTING & ENGINEERING**

### SUBSURFACE EXPLORATION

### LITHOLOGIC LOG OF DB-19

Client · French Limited Task Group
Project Name : French Limited Design Phase
Project Location : Crosby, Texas
Job Number : 2870—017 Bonng Na : D
Logged By · Cameron Teller
Approved By : Hoyt Clark
Drilled By : Fugro Geosciences, Inc.

Bonng Na : DB-19

DRILLING AND SAMPLING INFORMATION

Date Started: 9-13-90 Date Completed: 9-14-90
Method: Mud Rotary Total Depth 60 ft
WELL COMPLETION INFORMATION
Screen Dia: Length:
Slot Size: Type:

Casing Dia: Length:

	Cosing Did :										
DEPTH N FEET	DESCRIPTION	SAMPLE NO	SAMPLE	1000	ENICAL ALYSIS	OTECH ALYSIS	CRAPHIC LOC	BORING COMPLETION			
2 %	SURFACE ELEVATION: 14,5	SANI	35		5	ಆ₹	8	<b>8</b> €	1		
	Fill; stiff dark brown silty clay, same sand, roots.										
		1	ST		7		$\blacksquare$	1	ì		
5 —	5 9.5	2	ST				曲	1	1		
	Loose brown medium sand, little gravel, little silt.	3	ST	++			••		- 1		
	Soft brown clayey silt.  Very stiff dark brown to black sandy clay.		<del> }</del>		4		呼	1	ı		
10	CNC	4	ST		4				4		
		,-									
	Loose brown fine to medium sand, some silt	5	ss	-   -	7		4	1	- [		
15			33		~	$\dashv$			+		
3					ļ						
=		6	SS		7						
20 -	5			11	$\dashv$				4		
1								ļ	į		
=		7	SS		7			1	1		
25 -							-		1		
=	<u>28</u>				_						
30 <del>-</del>	Firm gray sity day	8	ST		1			}			
~ =	28 -13.5 V Firm gray sity day  C1  -19.5 V	9	ST	_	1	1		u l	7		
3	-19,5 V	10	ST		7			JECK	ı		
35 🚽	Dense tan silt.	11	ST		7	7		CROUT TO SUPFACE	4		
#	-gray below 36 ft, moderate odor	12	ST					됩	ł		
=		13	-	1-1	-	$\dashv$		8	ı		
40-	Medium dense gray fine sand and sit.		<del></del>	<del></del>	4	$\dashv$	₩₩	İ	$\exists$		
3	- T	14	$\rightarrow$		4		Ш				
#	1MI	15	ST		4	_	1111		- 1		
45-		16	ST	1.	4		Ш		+		
4	Note: #100 is a duplicate of #15	17	ST	- [ - [.	1		Щ				
]	-some clay below 48 ft	18	ST		1		Ш	Ì	ł		
50-]		19	ST		7		$\parallel \parallel$		1		
=	1	<u> </u>	ST	++	7	$\dashv$					
55 –			<del>- i</del>		+	{	Ш				
35	$\mathcal{M}$		ST		4		111	-	1		
ᅾ	58 - 43.5 V	22	ST		4		Щ				
Stiff rea-brown, yellow, gray silty clay, laminated, horizontal fine sand 23 ST and silt partings, variable plasticity.											
SAMPLER TYPE  SS - DRIVEN SPUT SPOON RC - ROCK CORE  C2 HSA - HOLLOW STEM AUGER DC - DRIVING CASING											
	RESSED SHELBY TUBE CT - CONTINUOUS TUBE CFA - C	CONTINUOU	S FUG	HT AUGERS				ID DRILLIN			



COMPANY : FUGRO GEOSCIENCES INC. OTHER SERVICES. HELL : DB-19 LOCATION/FIELD : FRENCH LTD. COUNTY : HARRIS STATE : TEXAS SECTION TOUNSHIP RANGE : : 09/14/90 PERMANENT DATUM : GL ELEVATIONS DATE DEPTH DRILLER : 60 J.M. ELEV. PERM. DATUM: GL KB : GL LOG BOTTOM 59.60 LOG MEASURED FROM: GL DF : GL LOG TOP -5.00 DRL MEASURED FROM: GL GL : GL CASING DRILLER : 00 LOGGING UNIT : 9002 CASING TYPE : 00 FIELD OFFICE : TULSA.OK. CASING THICKNESS. 00 RECORDED BY : 1.F.ALANIZ BIT SIZE : 3.874 BOREHOLE FLUID : H2O/GEL FILE : ORIGINAL MAGNETIC DECL. : 5.8 RH : 0 TYPE : 9030AA MATPIX DENSITY RM TEMPERATURE : 271 : 0 LOG MATRIX DELTA T PLOT : COTTER 0 FLUID DENSITY : 1.1 : 57

### ALL SERVICES PROVIDED SUBJECT TO STANDARD TERMS AND CONDITIONS

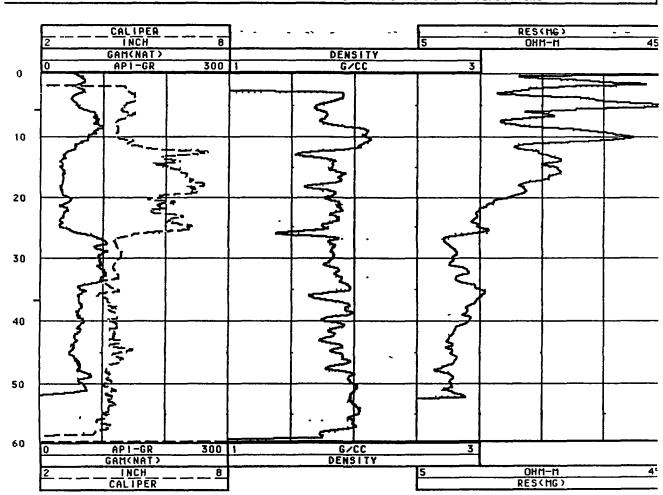
FLUID DELTA T

: 189

THRESH: 50000

NEUTRON MATRIX : SANDSTONE

REMHRKS



## minut -

٠,

### Appendix 5

Boring Log and Well Construction Details for REI-6-1

01502437



### SUBSURFACE EXPLORATION **RECORD**

	FRENCH LIMITED TASK FORCE								REI:6-	1				
	Archisect Engineer C. Itin						_ Je		27502					
(	Proper Name French Site			Drewn By JB										
(	Project Location Crosby Texas						_ 🗛	-	v_JDA_					
	DRILLING and SAMPLING INFORMATION Sunset Will 1	2N 40	•	-					TES	T DATA				
		30		n.										
1	Dritt Foremen_G_LittelBreen Sempler DI	2		•				E	y_*					
(	Rock Core Dis.			n.				<b>§</b>		•	<b>₽</b>		<b>1</b>	
(	Bering Method RW/HSA Bretty Tube DD		۹,	¥	ENY	GROUND WATER		1 1 Com	AMP. Cray bea.	Netwood Ony Demark Pot./Eu ft.	Consent			
	SOIL CLASSIFICATION	Į.	DEPTH	SAMPLE NO.	SAMPLE TYPE	S RECOVERY	OUND	Standard Per N, Blows/Ft.	Uncontra Strength Postal Po qp Tom/	- T	PA.A	Š.	3 6 8	
	SURFACE ELEVATION - 13.0	200	28	35	3	*	Š	8 Z	43 BE	_01 X	Ž	ALPRA	コペス	
П	SILTY CLAYEY SAND, fine grained, brown to gray, with some thin silty clay seams.		=	01	ss	39		10					E	
耳	(SH-SC) $(-)200 = 202$			02	ss	56		8					<b> </b> =	
=	· (-)200 = 12.6%		=											
I		4.5	=	03	SS	56		4					=	
H	SAND, medium to fine grained,		5-	nμ	SS	67		3					=	
	gray.	}	=					1				ŀ		
兰	(-)200 = 5.2X (-)200 = 2.9X		_		-	_		_			•	]	F	
=	( )=00 = 21,72			06	SS	67		7				•	l F	
$\exists$			10-	07	SS	5.4		11				Í	I E	
П			=					**			1		l E	
$\exists$			-	08	SS	39		15				1	1 E	
긬	(-)200 <b>-</b> 2.9%		-	09	ss	33		18					<del> </del>	
	(-)200 = 2.6X	1	=	1	-	-		١,,	l		l	1	1 F	
	(SP-SW) (-)200 = 1.8%		15_	<del> </del>	SS	56		16			1			
1111	(-)200 <b>-</b> 2.8%		=	11	SS	61		11	l		1	}	1 F	
_=	(-)200 - 2.1	Į.	=	12	ss	67		8			l	<b>[</b>	IE	
$\Box$	(-)200 <b>-</b> 2.4%	1	·=	}	133	<del> </del>	•	ľ	Ī	1	[	1	IE	
$\exists$		I	]	13	SS	56	]	10	1		1		I É	
日		ł	20-	14	SS	61		12		]			=	
111				1	ss	_		12					F	
	SILTY CLAY, yellow brown to		] =	16	ss	67	1	13			ł		IE	
	gray mottles (CH-CL)	25.0	25	<del>  -  </del>	1-	<del>  -</del>	1							
=			43-	17	SS	56	_	5	<u> </u>	<u></u>		<u> </u>	1	
Ξ	Continued on page 2			1							1		1 =	
			•	1		1		<u> </u>	1	1	<u> </u>	1	<u></u>	

SAMPLER TYPE SS - DRIVEN SPLIT SPOON
ST - PRESSED SHELBY TUBE
CA - CONTINUOUS FLIGHT AUGER
RC - ROCK CORE

GROUND WATER DEPTH T AT COMPLETION T AFTER FT WATER ON RODS

BORING METHOD MUNING METHOD

HSA - HOLLOW STEM AUGERS

CFA - CONTINUOUS FLIGHT AUGERS

DC - DRIVING CASING

MD - MUD DRILLING



page 2 of 3

### SUBSURFACE EXPLORATION **RECORD**

	Chem FRENCH I IMITED TASK FORCE						. 6	pring - F	REI:6-	1	<del></del>			
	Arehisset Engineer C. It1n				Job - 27502									
	Proper Name French Site				Drown By. UD									
	Project Lection_Crosby Texas						_ 4	pprovoq 3						
,	DRILLING and SAMPLING INFORMATION  Date Standon 3-2-84 Hammer Wit.	9N 140		••					TES	T DATA				
	Dete Completed 3-3-84 Hammer Dren	30	······································	. [										
	Dreft Foreman G. Littel Secon Servetor Of	2						Ŧ		1				
	Inspector IR Rock Core Dis.	_		R.				Ē					=	
	Boring Method RW Shelby Tube OD_	3	YPE	ENV	GROUND WATER		Tour Control	SHETY CITY SHE	Domety P.	Į.	Limit Limit 1998 Len			
	SOIL CLASSIFICATION	\$ e	11 12	SAMPLE NO.	SAMPLE TYPE	S RECOVERY	045	Standard Per N, Bloss/Ft.	Unconinud Strongth qui Pocho: Program qj? Tomi/Ft.	7	Network Dry No /cu	r Consent	7 T. 20 J. 2	
	SURFACE ELEVATION —	S C	DEPTH SCALE	₹2.	248	<b>3.</b> R	ě	% ¥.	Si ke	7 to	į		35%	
11	SILTY CLAY, yellow brown to gray,		1		1			_	į l			ا۔۔ ۔۔ا	LL=70	
1	mottles, with thin silt & sand		11	18	SS	94		7			1	36.32	PI-57	
	ch-cl) (-)200 = 98.1%		1	19	SS	67		21	]		1	35.7%		
-	(-)200 = 81.3		_		_				i	•	ļ		PL-21 -	
		29.7	30-	20	SS	33		11	1		ł		PI-28	
	SANDY CLAYEY SILT .		30-3	21	SS	67		27		'		ĺ	F	
	very fine grained, olive to			-		-			] [		•	I	l F	
	gray.			22	SS	67	. 1	31	1 1		l	1	l E	
1111	(HL) (-)200 = 54X			23	SS	67		20		•	ļ		l F	
	(-)200 <b>-</b> 57X		35		-	<u>:</u>					Į	1	l F	
	/ 2000 - 40 <b>7</b>			24	SS	67		23			}	Į.	l F	
7	(-)200 <b>= 49</b> X	37.5	_	25	SS	89		23			1	1	l F	
7	THE MAN AND STATE OF THE PARTY	37.3		23	33	07		23			1	l	I E	
	SILTY CLAYEY SAND, fine grained, very fine grained, olive gray, with thin silty clay seams, a			26	SS	78		22	Į,		l	l	F	
	with thin silty clay seams, & oily streaks and stains.	ľ	40_	27	SS	22		20	ł		l	ł	l F	
1		ŀ	-	21	33	23		20			1	l		
111	(SC-SH) (-)200 = 53% (-)200 = 41%	<u> </u>	·.=	28	ss	67		27	ì	Į .	1	1		
二	(-)200 = 412 (-)200 = 60X	<b>l</b> '	=	-	1	<u> </u>			1		į	1		
=	. /		=	24	SS	67		28	I	•				
I	0 to 26 40°		_ =	30	ss	67		48	1	1		1		
ㅋ	JULY BUK STAGKSDY'L	746	45—	<del>                                     </del>	<del> </del>	<del> </del>			ļ.	1	1	1	=	
=	VERY SILTY SANDY CLAY/CLAYEY	1	=	31	SS	83		37	1	1	]	20.8	L-26	
	SANDY SILT. red brown to gray			32	ss	89		20	1	1	1		PL-18	
I	with thin silt & sand seams. (CL-ML) (-)200 = 84%		=	_	1-				1	ł		1	P1-8 E	
	(-)200 = 51%		=	33	ss	95	1	17	1	}	}	1	1-23	
	Change to Silty Clay	51.0	50	34	SS	95		19	1	1		16.9	PL=15	
1	Continued on page 3		-	1	1		1				1	1	1 E	
		<u> </u>		1	<u></u>		1_	<u></u>		1	<u> </u>			

SAMPLER TYPE

SE — DRIVEN SPLIT SPOON

ST — PRESSED SHELBY TUBE

CA — CONTINUOUS FLIGHT AUGER

RC — ROCK CORE

V AT COMPLETION HRS

WATER ON RODS

GROUND WATER DEPTH 5.4 FT.

FT. FT.

BORING METHOD

MSA - MOLLOW STEM AUGERS
CFA - CONTINUOUS FLIGHT AUGERS
DC - DRIVING CASING
MD - MUD DRILLING



### SUBSURFACE EXPLORATION **RECORD**

•	French Limited Task Force						_ 84	rine o	REI:	6-1				
Anghinget Engineer C. Tr.Ltp.				٠,		275-	-02							
French Site					_ D	nen Bv.	JB							
Crosby, Texas					_ &	proved B	- JDA							
-	DRILLING one SAMPLING INFORMATIO	M								T DATA				
	3-3-85 Harring Wi.	14		<b>bu</b> . ,		-	_							
ŧ	Sale Completed	3		۹.						~.				l
•	Ordi Foreman G. Littel Boson Someter OD		<u>2</u> ,	۹.				Ę	¥. :					ĺ
•	Rect Care Da	_	-	n.				•	E .	1	<b>2</b>		ī	ĺ
•	forms MethodShifty Take OD_		30	٨.		•				2	2	_	117	İ
	ر المرابع المرابع المرابع المرابع المرابع المرابع المرابع المرابع المرابع المرابع المرابع المرابع المرابع المر				Ε	E.	8	12		15	2 2	Constitution	433	ł
- 1	SOIL CLASSIFICATION	<b>F</b> .	2	2	3	Š	暑	11		1	ed Ory Be Are.	8.	1 3 5	
ł	SURFACE ELEVATION -	100	SCAL	SALPPLE NO.	SAMPLE TYPE	& RECOVERY	GROUND WATER	Z Z	20 20	× 0	1	Photo	7.0	ĺ
				3 X	3	*	ō	& z	34 89	Z×.	2	2	コイス	
コ	SILTY SANDY CLAY/CLAYEY SILT	كىلگ	=						f :					
$\exists$	(CL-HL)		1 =	35	SS	95		16					LL=62	F
ㅋ	SILTY CLAY, redbrown to gray		-	7		~	1		•				PL=21	F
=	mottles, stiff with silt seams		=	30	SS	33		19					PI-41	E
Ħ	(-)200 <del>-</del> 987		] =	37	ss	95		17	•	•			, ,	E
님	(CH-CL)		55—		-		1	• •	ł .				LL-39	F
7	(on out	56.5	] =	38	SS	95		20			į	ł	PL-20	F
$\exists$	VERY CLAYEY SILT, redbrown to		1 =			_	1	_			1	ì	PI-19	
コ	gray, stiff with very thin			39	SS	100	1	9	<b>!</b>		1	1	i	E
$\exists$	silt seams	1	=	40	ss	O.E.	1		ĺ	i .	1	[	LL-51	F
コ	(HH)		-		33	77	1	9	Ì	Ì	ì	ł	PL-20	
듸	(im)		60_	41	SS	75	ı	9	i	ł	ì	Ì	PI=19	F
J			1 =	<u> </u>		-			1				<del> </del>	t
3	Boring Terminated at 61ft.		= ا	•						l	1	1	1	E
4	(EI48.0)	i	-	•				i		1	[	•	i	
=	•		=	1		Ì			į	ł	l	1	1	F
コ		ł	65-	1		ł			ł	I	}	1	1	L
$\exists$	1 ( 0(1)	}	)" <u>=</u>	}	•	1			}	j	)	}	3	E
4	11 50600	ł	=	ł	İ				l l		•	1		F
듸	1/60C. = 3/31	[	_	1	•				l .	•	1	1	1	F
コ	WERL. 13 SCEEN 6-1 25 578 6-2	ł	[ =	j	ł	ł	l		ł	l	ł	i		t
-	6-1 25 5-00		-	1	1	1	1	ŀ	1	Į	1	1	1	H
コ	6-7		70-	1	1	]	}	}	}	ł	j .	1	1	F
Ⅎ	<b>O</b> -	ł	l :	1	ŀ	1	1	1		I	1	ł	1	t
4			! :	ł	ſ	l	ı	Ì	ł	ł	ł	ł	i	Ł
コ			-	7		ł	1	ł	1	ł	1	ł	1	F
ᆿ			75	1			1		}	l .	}		1	F
크	•	}	=	j	1	1	1	1	1	1	l .	1	1	t
耳		1	75_	3	1			1	١.	{	1	1		F
111/111		i	:	4	1	1	1	I	1:	1	1	1	1	F
=		l	1:	1	i	1	1	ł	1	1	1	1	1	t
		<u> </u>		1	<b>S</b>	1	1	<u> </u>	<u>. f</u>	1	1	1		_

SAMPLER TYPE
SS — DRIVEN SPLIT SPOON
ST — PRESSED SHELBY TUBE
CA — CONTINUOUS FLIGHT AUGER
BC — BCCY CORE

GROUND WATER DEPTH

3.4 T AT COMPLETION T AFTER HRS WATER ON BOOK

FT. FT

BORING METHOD

HEA - MOLLOW STEM AUGERS

CFA - CONTINUOUS FLIGHT AUGERS

DC - DRIVING CASING

MD - MUD DRILLING

DETAILS OF MONITOR WELL CONSTRUCT		
Project Name: FRENCH LIMITED SITE Boring Number: Project Number: 275-02 Date Installed		
Water Level Measurement: 6.85		
Top of Casing Elev 14.78		
Water Level Elev.	•	
Protective Steel Cover		
2.35 2.84		••
Ground Surfa		12.43 DESCRIPTION
/ // ES/	1	STITY CLAVEY
8" inch diameter Surface Cas	Ing 4.5	SAND (SC)
	į	571 7V 54115
Cement Bentonite Grout (4-1 mix)	-	SILTY SAND fine to medium
		grained with
51		some gravel
		(SP-SW)
27' - depth of Surface Casing	25.0_	-12.6
		SILTY CLAY stiff
4" inch PVC Well Casing		.= :
1 NN - CI	2	(CH)
27 top of seal		
29 - bottom of seal  30 = top of screen		
29 - bottom of seal	_ 29.7	
30 = top of screen		SANDY SILT very fine
	ļ	, and the second
		(ML)
Sand Pack		
Slotted Well Screen /NT	37.5	CTITU CLAVEY
		SILTY CLAYEY SAND, with some
		thin clay seams
		(SC-SM)
	45.0	-33.6
50 - bottom of screen C	2	VERY SILTY SANDY
51 - Total Dept		CLAY (CL-ML)
8"- Borehole Diameter		SILTY CLAY
		stiff (CH)
L	urce ent	GINEERING

# BOOKMARK

## Appendix 6

Analytical Reports
INT-11B and INT-11C Soil samples

01502438

Report Prepared for

FLTG, INC.

1024 GULF PUMP ROAD. CROSBY TX 77532

Attention : TED DAVIS

by

Keystone Lab - Houston
A Division of CHESTER LabNet
8300 Westpark Drive
Houston, Texas 77063
(713) 266-6800

CERTIFIED BY: Domit D'Ostolanie

Dan Pastalaniec Project Manager

> WORK ORDER : H92-04.243 DATE RECEIVED : 22-APR-1992

> > In 1/12]

R 6-5-00

'ROJECT ID :
'.O. NUMBER : SPEC0018

FRENCH LTD. PROJECT

FLTG, Incorporated 15010 FM 2100, Suite 200 Crosby, Texas 77532

(713) 328-5860 FAX (713) 328-2996

DATA PACKAGE TO: FLTG, Incorporated Crosby, TX 77532 1024 Gulf Pump Road REPORTING LABORATORY

Name: Keystone Lab-Houston

Address: 8300 Westpark

Houston, TX 77063

Contact: Daniel D. Pastalaniec

Phone: (713) 266-6800 (713) 974-5491 FAX:

Requested By: TD

Standard TA: Y

Days: 21

Notified by:

Date:

FLTG MATRIX CODE: SPEC FLTG SET NUMBER: SPECO018 Quality Control Level: I

Site Location: 1024 Gulf Pump Road Crosby, TX 77532

FLTG Sample No	Date	Time	Location	Grab/ Comp	Matrix	Туре
SPEC0018-01 SPEC0018-02	4/21/92 4/21/92	15:30	INT-11B INT-11C	Grab Grab	Soil Soil	ENV ENV
			-			
		-				

SAMPLER'S Name(s) (PRINT) 1 - Donald A. Flory 2.

Affiliation Drilling Crew

Carrier:

з.

Bill No.:

RELINQUISHED BY

(SIGN) (INITIALS) Date Time

(INITIALS) (SIGN)

Date Time

RETURNED BY

Time

(SIGN) (INITIALS) Date

DISPOSED BY

RECEIVED BY

(SIGN) (INITIALS) Date lime

1. 2.

NOTES TO LAB:

Run naphthalene in VOA

( milar) (White - Original Yollow - Sot D.

No.: 277 Page 2 of 2

	FLTG MATRIX CODE:SPEC Lab Work Order No: FLTG SET NUMBER:SPEC0018								
SAMPLING SOP	REFERENCE:N/A								
DESCRIPTION:	DESCRIPTION:Soil boring								
NOTES TO SAME	NOTES TO SAMPLER:								
	Pr	ARAMETER	S REQUE	STED					
GC8015 (TF	<b>ЭН) Т</b> О	DC		TOX		VOA-SX (NAF	PHT )		
FLTG Sample No	Lab Sample ID	Conta	iner			Preserva	ative		
SPEC0018-01 SPEC0018-02			M glass M glass			Refrig Refrig			
							;		
COMMENTS			(SI	et QC C GN)	Complet (INIT	ed by: IALS)	Date		

# KEYSTONE LAB - HOUSTON A Division of CHESTER LabNet

Reported on: 8-MAY-1992

Client Name: FRENCH LIMITED Work Order: H92-04.243
Sample ID: HAT0427-019 Date Collected: 21-APR-1992

Sample Name : H92-04.243-002 Matrix : SOIL

Project No.: SPECO018

Date Received: 22-APR-1992

Percent Moisture: 19 %

Date Extracted: 23-APR-1992

Client ID : SPECO018-01 INT-11B Checked by :

Organic Analysis Data Sheet Compounds Analysis by SW846 Method 8015

Date Analyzed: 27-APR-1992 20:26 Dilution Factor: 1.000

Analyzed by : 8388624

Detection Detected
Cas # Compound Limits Conc. mg/kg

TOTAL PETROLEUM HYDROCARBONS 620 97100

U = Undetected at the Listed Detection Limit .

J = Compound is present, but below the Detection Limit.

B = Compound is also found in Blank.

# KEYSTONE LAB - HOUSTON A Division of CHESTER LabNet

Reported on: 8-MAY-1992

Client Name: FRENCH LIMITED Work Order: H92-04.243
Sample ID: HAT0427-022 Date Collected: 21-APR-1992

Sample Name: H92-04.243-005 Matrix : SOIL

Project No.: SPECO018 Date Received: 22-APR-1992
Percent Moisture: 19 % Date Extracted: 23-APR-1992

Client ID : SPECO018-02 INT-11C Checked by :

Organic Analysis Data Sheet Compounds Analysis by SW846 Method 8015

Date Analyzed: 27-APR-1992 22:05 Dilution Factor: 1.000

Analyzed by : 8388624

Detection Detected
Cas # Compound Limits Conc. mg/kg

TOTAL PETROLEUM HYDROCARBONS 620 69500

U = Undetected at the Listed Detection Limit .

- J = Compound is present, but below the Detection Limit.

B = Compound is also found in Blank.

Page

#### Summary of Analytical Results

Date received: 22-APR-1992 Job name: H92-04.243 Customer: FLIG, INC.

		Samples					
Keystone ID Sampling Point Date Sampled Customer ID Parameters		243-001 QA QC 21-APR-1992 LAB BLANK	243-002 X 21-APR-1992 SPE00018-01 NA INT-118	243-003 QA QC 21-APR-1992 SPEC0018-01 DUP (INT-11B dup analysis)	243-004 QA_QC 21-APR-1992 SPEC0018-01 MS	243-005 X 21-APR-1992 SPEC0018-02 NA INT-11C	
Total Organic Carbon (Solid) Analyst: ELL Date/Time: 5/01/92 11:31 Dilution: 20	ng/Kg	, <50.0	38900	39200	105*	31700	
Notal Organic Halogens (Solid) Analyst: RNS Date/Time: 5/05/92 08:30	ng/Kg	<10.0	80300	79000	101#	112000	

Dilution: 5000

<sup>\* - %</sup> Recovery NR - Not Required

NA - Not Applicable # - Blank Spike Recovery

KEYSTONE DC # ---- 8

SAMPLE NUMBER: SPECO018-01 INT-

#### ORGANIC ANALYSIS DATA SHEET -

ABORATORY NAME: KEYSTONE ENV.

AB SAMPLE ID NO.: 910424302

CASE NO. : ----

INITIAL CALIBRATION DATE: ......

CONTRACT NO.: --

8 .

...... 250

DATE SAMPLE RECEIVED: 04/22/92

**YOLATILES** 

ONCENTRATION: LOW DATAFILE: 4U04243V02 ATE ANALYZED: 05/01/92 DILUTION FACTOR: 1.60

DETECTION AD 5//92 AMOUNT LIMITMILL COMPOUND FOUND (MICROGRAMS / KG) CO10 CHLOROMETHANE CO15 BROMOMETHANE 16 U 16 U CO20 VINYL CHLORIDE 16 CO25 CHLOROETHANE 16 CO30 METHYLENE CHLORIDE 8 CO35 ACETONE 16 U CO40 CARBON DISULFIDE 8 CO45 1,1-DICHLOROETHENE 8 300 COSO 1,1-DICHLORDETHANE 7200 CO53 1.2-DICHLOROETHENE (TOTAL) 8 CO60 CHLOROFORM 8 ...... 12000 8 CO65 1, 2-DICHLORDETHANE 16 U C110 2-BUTANONE вυ C115 1, 1, 1-TRICHLOROETHANE 8 ..... 7000 C120 CARBON TETRACHLORIDE C125 VINYL ACETATE 16 U C130 BROMODICHLOROMETHANE 8 U 8 U C140 1.2-DICHLOROPROPANE C143 CIS-1, 3-DICHLOROPROPENE 8 U C150 8 ..... 700 TRICHLOROETHENE 8 U C155 DIBROMOCHLOROMETHANE C160 1,1,2-TRICHLORDETHANE 12 C165 BENZENE 27 8 U C172 TRANS-1, 3-DICHLOROPROPENE C175 2-CHLOROETHYLVINYLETHER 16 U C180 BROMOFORM 8 U C205 54 4-METHYL-2-PENTANONE 16 C210 2-HEXANONE 16 7 J C220 TETRACHLOROETHENE 8 ......... 3700 1, 1, 2, 2-TETRACHLOROETHANE 8 130 C225 8 ...... C230 TOLUENE 8 U C235 CHLOROBENZENE 8 ..... C240 ETHYLBENZENE 8 U C245 STYRENE

XYLENES (TOTAL)

C250

J = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

KEYSTONE DC # ---- B

SAMPLE NUMBER: SPECO018-01 I

ORGANICS ANALYSIS DATA SHEET - PAGE 4

LABORATORY NAME: KEYSTONE ENVIRONMENTAL RESOURCES, INC. CASE NO.: ----

3C REPORT NO.: .....

ANALYST: PVG

DATAFILE: 4004243V02

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS # VOLATILE COMPOUND NAMES

PURITY AMOUNT SCAN#

76-01-7 ETHANE, PENTACHLORO-

738 891 690

UG/L

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

KEYSTONE DC # ---- 8 LF NUMBER: SPECOO18-01 INT-11

SAMPLE NUMBER: SPECO018-01 INT-118

#### ORGANICS ANALYSIS DATA SHEET

LABORATORY NAME: KEYSTONE ENV.

LAB SAMPLE ID NO.: 910424302

SAMPLE MATRIX: WATER

DATA RELEASE AUTHORIZED BY: . /

CASE NO. : ----

INITIAL CALIBRATION DATE: 04/30/92

CONTRACT NO.: --

DATE SAMPLE RECEIVED: 04/22/92

YOUATILES

CONCENTRATION: LOW

DATE ANALYZED: 05/04/92

DATAFILE: 4U04243V02C

DILUTION FACTOR: 310.00

		COMPOUND	DETECTION LIMITMILL; CMICROCRAMS	AMOUNT  KAN FOUND FU 5/4/52  / LITER )
	CO10	CHLDROMETHANE BROMOMETHANE VINYL CHLORIDE CHLOROETHANE METHYLENE CHLORIDE ACETONE	3100 U	سے نہیے میں بیپار ''ناہ میت ہاللہ 'کا 'جی باک کا اسا گاہ گا۔ کا انسان ہالہ
		BROMOMETHANE	3100 U	
	C020	VINYL CHLORIDE	3100 U	
	CO25	CHLOROETHANE	3100 U	
	CO30	METHYLENE CHLORIDE	1550 U	
	C035	ACETONE CARBON DISULFIDE  1.1-DICHLOROETHENE 1.1-DICHLOROETHANE 1.2-DICHLOROETHENE (TOTAL) CHIOROEDRM	3100 U	
	C040	CARBON DISULFIDE	1550 U	
	CO45	1,1-DICHLOROETHENE	1550 U	
	C050	1,1-DICHLOROETHANE	1550	520 J
	C053	1,2-DICHLOROETHENE (TOTAL)	1550	5000
	C040	CHLOROFORM	1550	27000
	C065	1.2-DICHLOROETHANE	1550	21000
	C110	2-BUTANONE	3100 U	•
	C115	CARBON DISULFIDE  1, 1-DICHLOROETHENE  1, 1-DICHLOROETHANE  1, 2-DICHLOROETHANE  CHLOROFORM  1, 2-DICHLOROETHANE  2-BUTANONE  1, 1, 1-TRICHLOROETHANE  CARBON TETRACHLORIDE  VINYL ACETATE  BROMODICHLOROMETHANE  1, 2-DICHLOROPROPANE  CIS-1, 3-DICHLOROPROPENE  TRICHLOROETHENE  DIBROMOCHLOROMETHANE  1, 1, 2-TRICHLOROETHANE  BENZENE  TRANS-1, 3-DICHLOROPROPENE  BROMOFORM	1550 U	
	C120	CARBON TETRACHLORIDE	1550	26000
	C125	VINYL ACETATE	3100 U	
	C130	BROMODICHLOROMETHANE	1550 U	
	C140	1,2-DICHLOROPROPANE	1550 U	
	C143	CIS-1.3-DICHLOROPROPENE	1550 U	
	C150	TRICHLOROETHENE	1550	700 J
	C155	DIBROMOCHLOROMETHANE	1550 U	
	C160	1, 1, 2-TRICHLOROETHANE	1550 U	
	C165	BENZENE	1550 U	
	C172	TRANS-1, 3-DICHLOROPROPENE	1550 U	
	C180	BROMOFORM	1550 U	
	C205	4-METHYI -2-PENTANINE	3100 U	
	C210	2-HEXANONE TETRACHLOROETHENE 1, 1, 2, 2-TETRACHLOROETHANE	3100 U	
	C550	TETRACHLOROETHENE	1550	13000
	C225	1, 1, 2, 2-TETRACHLORDETHANE	1550 U	
	C230	TOLUENE	1550 U	
_	C235	CHLOROBENZENE	1550 U	
	C240	ETHYLBENZENE	1550 U	
	C245	STYRENE	1550 U	
	C250	XYLENES (TOTAL)	1550 U	

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

**VST ID: 4000** 

KEYSTONE DC # ---- 8

SAMPLE NUMBER: SPECO018-01 I

ORGANICS ANALYSIS DATA SHEET - PAGE 4

ABORATORY NAME: KEYSTONE ENVIRONMENTAL RESOURCES, INC. CASE NO.: ----

C REPORT NO.: ..... ANALYST: PVG DATAFILE: 4U04243V020

B. TENTATIVELY IDENTIFIED COMPOUNDS

SCAN# PURITY AMOUNT

AS # VOLATILE COMPOUND NAMES

UG/L

NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT. STD.

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

SAMPLE NUMBER: SPECO018-01 INT-11B



#### ORGANICS ANALYSIS DATA SHEET

LABORATORY NAME, KEYSTONE

LAB SAMPLE ID NO.: 920424302

SAMPLE MATRIX: SOIL

DATE ANALYZED. 05/27/92

DATA RELEASE AUTHORIZED BY:.

C540 ACENAPHTHYLENE

C545 3-N] TROANILINE

C555 2,4-DINITROPHENOL C560 4-N1TROPHENOL

C570 2,4-DINITROTOLUENE

C543 2,6-DINITROTOLUENE

C550 ACENAPHTHENE

C565 DIBENZOFURAN

CASE NO. . --

QC REPORT NO.: 05/18/92

CONTRACT NO.: ---

DATE SAMPLE RECEIVED: ..

610 J

360 J

SEMIVOLATILES

CONCENTRATION: LOW DATE EXTRACTED . . . .

DATAFILE: 4U04243C02

10000 U 50000 U

10000 ..

50000 U

50000 U 10000 ....

10000 U

10000 U

		COMPOUND	DETECTION LIMIT	AMOUNT FOUND
		COIN COILD	(MILLIGRAMS	KG)
	C315		10000 U	
	C325	BIS(2-CHLOROETHYL)ETHER	10000 U	
	C330	2-CHLOROPHENOL	10000 U	
	C335	1,3-DICHLOROBENZENE	10000 U	
	C340	1,4-DICHLOROBENZENE	10000 U	
	C345	BENZYL ALCOHOL	10000 U	
	C350	1,2-DICHLOROBENZENE	10000 U	
V	C355	2-METHYLPHENOL	10000 U	
_	C357	2,2'-OXYBIS(1-CHLOROPROPANE)	10000 U	
	C365	4-METHYLPHENOL	10000 U	
	C370	N-NITROSODIPROPYLAMINE	10000 U	
	C375	HEXACHLOROETHANE	10000	5800 J -
	C410	NITROBENZENE	10000 U	
	C415	ISOPHORONE	10000 U	
	C420	2-NITROPHENOL	10000 U	- <i>=</i> -
	C425	2,4-DIMETHYLPHENOL	10000 U	
	C430	BENZOIC ACID	50000 U	
	C435	BIS(2-CHLOROETHOXY)METHANE	10000 U	
	C440	2,4-DICHLOROPHENOL	10000 U	
	C445	1, 2, 4-TRICHLOROBENZENE	10000 U	
	C450	NAPHTHALENE	10000	1600 J
	C455	4-CHLOROANILINE	10000 U	
	C460	HEXACHLOROBUTADIENE	10000	38000
	C465	P-CHLORO-M-CRESOL	10000 U	
	C470	2-METHYLNAPHTHALENE	10000	350 J
	C510	HEXACHLOROCYCLOPENTADIENE		
	C515	2, 4, 6-TRICHLOROPHENOL	10000 U	
	C520	2, 4, 5-TRICHLOROPHENOL	50000 U	
	C525	2-CHLORONAPHTHALENE	10000 U	
	C530	2-NITROANILINE	50000 U	
	C535	DIMETHYL PHTHALATE	10000 U	

#### KEYSTONE DC # --- 8

SAMPLE NUMBER: SPECO018-01 INT-118

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

DATAFILE: 4U04243C02

DETECTION AMOUNT COMPOUND LIMIT FOUND (MILLIGRAMS / KG)	
C580 DIETHYL PHTHALATE C585 4-CHLOROPHENYL PHENYL ETHER C595 4-CHLOROPHENYL PHENYL ETHER C595 FLUORENE C595 4-NITROANILINE C610 4.6-DINITRO-2-METHYLPHENDL C615 N-NITROSDIPHENYLAMINE C625 4-BROMOPHENYL PHENYL ETHER C625 4-BROMOPHENYL PHENYL ETHER C630 HEXACHLOROPHENDL C630 HEXACHLOROPHENDL C635 PENTACHLOROPHENDL C645 ANTHRACENE C645 ANTHRACENE C645 ANTHRACENE C650 DI-N-BUTYL PHTHALATE C650 DI-N-BUTYL PHTHALATE C650 DI-N-BUTYL PHTHALATE C720 BUTYL BENZYL PHTHALATE C720 BUTYL BENZYL PHTHALATE C720 BUTYL BENZYL PHTHALATE C730 BENZO (A) ANTHRACENE C740 CHRYSENE C740 CHRYSENE C750 DI-N-DCTYL PHTHALATE C760 DI-N-DCTYL PHTHALATE C770 BENZO (A) PYRENE C770 BENZO (A) PYRENE C770 DENZO (A) PYRENE C770 DENZO (A) PYRENE C770 DENZO (A) H) ANTHRACENE C770 DENZO (A) H) ANTHRACENE C770 DENZO (CA, H) ANTHRACENE C	

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

SAMPLE: SPECOO18-01 INT-118

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME KEYSTONE ENV. RESOURCES

CASE NO.: --

QC REPORT NO.: 05/18/92 ANALYST: SC

DATAFILE: 4U04243CO

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS #	SEMIVOLATILE COMPOUND NAMES	SCAN#	PURITY	AMOUNT
				MG/KG
55880-7	7-8 1.3-BUTADIENE, PENTACHLORO-	604	900	5900
	ALIPHATIC HYDROCARBON	998		7300

ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

#### KEYS1 ONE DC # ---- B

SAMPLE NUMBER: SPECO018-02 IN

#### ORGANIC ANALYSIS DATA SHEET -

LABORATORY NAME: KEYSTONE ENV.

LAB SAMPLE ID NO.: 910424305

SAMPLE MATRIX: SOIL

CASE NO. : ----

INITIAL CALIBRATION DATE: ......

CONTRACT NO. : --

DATE SAMPLE RECEIVED: 04/22/92

VOLATILES

CONCENTRATION: **DATAFILE: 4U04243V05** LOW DATE ANALYZED: 05/01/92 DILUTION FACTOR: 1.60

	COMPOUND  METHANE	DETECTION & CMICROGRAM	ارامی AMOU FOU S / KG)	IND
CO10 CHLORO		16 U		
CO15 BROMOM	ETHANE	16 U		
CO20 VINYL	CHLORIDE	16 U		
CO25 CHLORO	ETHANE	16	5 J	
CO30 METHYLI	ENE CHLORIDE	<b>8</b>	. 49	
CO35 ACETONI	ENE CHECKIDE  DISULFIDE  CHLOROETHENE  CHLOROETHENE (TOTAL)  FORM  CHLOROETHANE  NONE  TRICHLOROETHANE  ACETATE  ICHLOROMETHANE  CHLOROMETHANE  CHLOROPROPANE  3-DICHLOROPROPENE	1 <u>6</u> U		_
CO40 CARBON	DISULFIDE	<b>8</b>	<u>3</u> J	l
CO45 1,1-DI	CHLOROETHENE	<u>8</u>	50	
CO50 1,1-DI	CHLOROETHANE	<u>8</u>	. 380	
CO53 1,2-DI	CHLOROETHENE (TOTAL)	<b>8</b>	. 2900	
CO60 CHLORO	FORM	8	8700	
CO65 1.2-DI	CHLOROETHANE	8	9100	
C110 2-BUTA	NONE	16	. 18	_
C115 1,1,1-	TRICHLOROETHANE	- 8	4 J	•
C120 CARBON	TETRACHLORIDE	· 8	. 5600	
C125 VINYL	ACETATE	16 U		
C130 BROMOD	ICHLOROMETHANE	8 U		
C140 1.2-DI	CHLOROPROPANE	- 8 U		
C143 CIS-1,	CHLOROPROPANE 3-DICHLOROPROPENE OROETHENE OCHLOROMETHANE TRICHLOROETHANE E 1,3-DICHLOROPROPENE			
C150 TRICHL	OROETHENE	8	480	
C155 DIBROM	OCHLOROMETHANE	B U		
C160 1,1,2-	TRICHLOROETHANE	8	g J	
C165 BENZENI	E	8	21	
	1,3-DICHLOROPROPENE	8 0		
		16 U		
C180 BROMOF	ORM	ευ		
	YL-2-PENTANONE	16	. 40	
C210 2-HEXA	NONE	16	4 J	ļ
	HLOROETHENE	8	. 3500	
	2-TETRACHLOROETHANE	8	. 100	
C230 TOLUEN	E	8 U 16 8 8 8 U 8 U	. 51	
	BENZENE	8 ប		
C240 ETHYLB			. 54	
C245 STYREN		ខប		
C250 XYLENES	S (TOTAL)	8	. 170	

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

KEYSTONE DC # ---- 8

ORGANICS ANALYSIS DATA SHEET - PAGE 4

\_ABORATORY NAME: KEYSTONE ENVIRONMENTAL RESOURCES, INC. CASE NO.: ----

SAMPLE NUMBER: SPECO018-02 I

3C REPORT NO.: .....

ANALYST: PVG

DATAFILE: 4U04243V05

B. TENTATIVELY IDENTIFIED COMPOUNDS

VOLATILE COMPOUND NAMES PURITY AMOUNT SCAN#

UG/L

76-01-7 ETHANE, PENTACHLORO-

894 736

340

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

[NST ID: 4000

KEYSTONE DC # ---- 8
SAMPLE NUMBER: SPEC0018-02 INT-11C

#### ORGANICS ANALYSIS DATA SHEET

LABORATORY NAME: KEYSTONE ENV. LAB SAMPLE ID NO.: 910424305

SAMPLE MATRIX: WATER

)ATA RELEASE AUTHORIZED BY:. J

CASE NO. : ----

INITIAL CALIBRATION DATE: 04/30/92

CONTRACT NO.: --

DATE SAMPLE RECEIVED: 04/22/92

**VO**CATILES

CONCENTRATION: LOW DATAFILE. 4U04243V05A

DATAFILE. 4U04243V05A

DILUTION FACTOR: 310.00

	COMPOUND	DETECTION LIMIT MELLE <del>LMIGROG</del> RAMS	AMOUNT  KA FOUND 574/92  / LITER )
C010	CHLOROMETHANE BROMOMETHANE VINYL CHLORIDE CHLOROFTHANE	3100 U	
C015	BRUMUMETHANE	3100 U	
6050	VINYL CHLURIDE	3100 U	
9020	···		
C030	MEINYLENE CHLURIDE	1550 U 3100	0100 1
C035			2100 J
C040		1550 U	
CO45		1550 U	
0050	1.1-DICHLOROETHANE	1550 0	
	1, 2-DICHLOROETHENE (TOTAL)	1550 U 1550 U 1550 1550	9300
	CHLOROFORM	1550	9200
C110	1,2-DICHLOROETHANE 2-BUTANONE	3100 !!	7200
· - C115	1 1. 1. TO TOUL ODDETWANE		
0110	- CARBON TETRACHLORIDE - VINYL ACETATE	1550	7500
C125	UINVI ACETATE	3100 U	7300
C130	BROMODICHLOROMETHANE	1550 U	
C140	1, 2-DICHLOROPROPANE	1550 U	
C143	CIS-1,3-DICHLOROPROPENE	1550 U	
		1550 U	
	DIBROMOCHLOROMETHANE	1550 U	
	1, 1, 2-TRICHLORDETHANE	1550 U	
	BENZENE	1550 U	
	TRANS-1, 3-DICHLOROPROPENE		
C180		1550 U	
C205	4-METHYL-2-PENTANONE		
C210	2-HEXANONE	3100 U 3100 U	
C250	TETRACHLOROETHENE	1550	4500
C225	1, 1, 2, 2-TETRACHLOROETHANE	1550 U	
C230	TOLUENE	1550 U	
C235	CHLOROBENZENE	1550 U	
C240	ETHYLBENZENE	1550 U	
C245	STYRENE	1550 U	
C250	XYLENES (TOTAL)	1550 U	

<sup>1 =</sup> UNDETECTED AT THE LISTED DETECTION LIMIT

<sup>1 =</sup> COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

KEYSTONE DC # ---- 8

ABORATORY NAME: KEYSTONE ENVIRONMENTAL RESOURCES, INC. CASE NO.: ----

SAMPLE NUMBER: SPECO018-02 I

C REPORT NO.: .....

ANALYST: PVG

ORGANICS ANALYSIS DATA SHEET - PAGE 4

DATAFILE. 4U04243V05A

B. TENTATIVELY IDENTIFIED COMPOUNDS

VOLATILE COMPOUND NAMES

SCAN# PURITY AMOUNT

UG/L

NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT. STD.

J = ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED

SAMPLE NUMBER: SPECO018-02 INT-11C

ORGANICS ANALYSIS DATA SHEET -

LABORATORY NAME: KEYSTONE

LAB SAMPLE ID NO : 920424305

SAMPLE MATRIX: SOIL

DATA RELEASE AUTHORIZED BY:.

CASE NO.: --

QC REPORT NO.: 05/18/92

CONTRACT NO.: ---

DATE SAMPLE RECEIVED:

**SEMIVOLATILES** 

DATE EXTRACTED: .....04/23

C543 2,6-DINITROTOLUENE

DATE ANALYZED: 05/24/92

**DATAFILE: 4U04243C05** 

10000 U

			DETECTION	AMOUNT
		COMPOUND	LIMIT	FOUND
			/// Comm x 4/// // // // // // // // // // // // /	KG)
	C315	PHENOL	10000 U	به ۱۳۰ مده انتقا سه خبه مبه سه <sub>جب</sub> ه هم هم جب هم هم سه سال المث
	C325	BIS(2-CHLOROETHYL)ETHER	10000 U	
	C330	2-CHLOROPHENOL	10000 U	
	C335	2-CHLOROPHENOL  1,3-DICHLOROBENZENE  1,4-DICHLOROBENZENE  BENZYL ALCOHOL  1,2-DICHLOROBENZENE  2-METHYLPHENOL	10000 U	
	C340	1,4-DICHLOROBENZENE	10000 U	
	C345	BENZYL ALCOHOL	10000 U	
	C350	1.2-DICHLOROBENZENE	10000 U	
	C355	2-METHYLPHENOL	10000 U	
	C357	2,2'-OXYBIS(1-CHLOROPROPANE)	10000 0	
	C365	4-METHYLPHENOL	10000 U	
,	C370	N-NITROSODIPROPYLAMINE	10000 U	_
	C375	HEXACHLOROETHANE	10000	. 2200 J
	C410	NITROBENZENE	10000 U	
	C415	N-NITROSODIPROPYLAMINE HEXACHLOROETHANE NITROBENZENE ISOPHORONE 2-NITROPHENOL 2,4-DIMETHYLPHENOL BENZOIC ACID BIS(2-CHLOROETHOXY)METHANE	10000 U	
	C420	2-NITROPHENOL	10000 U	
	C425	2,4-DIMETHYLPHENOL	10000 U	
	C430	BENZOIC ACID	50000 U	
	C435	BIS(2-CHLOROETHOXY)METHANE	10000 U	
	C440	2, 4-DICHLOROPHENOL	10000 U	
	C445	1, 2, 4-TRICHLOROBENZENE	10000 U 10000	440
	C450	NAPHIHALENE	10000	. 660 J
	C455	4-CHLUKUANILINE	10000 U 10000	17000
	C460	2,4-DIMETHYLPHENOL BENZOIC ACID BIS(2-CHLOROETHOXY)METHANE 2,4-DICHLOROPHENOL 1,2,4-TRICHLOROBENZENE NAPHTHALENE 4-CHLOROANILINE HEXACHLOROBUTADIENE P-CHLORO-M-CRESOL 2-METHYLNAPHTHALENE	10000	. 17000
	C465	P-CHLUKU-M-CKESUL	10000 U	
	C510	HEXACHLOROCYCLOPENTADIENE	10000 U	
	C515 C520	2,4,6-TRICHLOROPHENOL 2,4,5-TRICHLOROPHENOL	50000 U	
	C525	2-CHLORONAPHTHALENE	10000 U	
	C530	2-NITROANILINE	50000 U	
	C535	DIMETHYL PHTHALATE		
_	C540	ACENAPHTHYLENE	10000 U	
	C545	3-NITROANILINE	50000 U	
	C550	ACENAPHTHENE	10000 U	
	C555		50000 U	
		2,4-DINITROPHENOL	50000 U	
	C560	4-NITROPHENOL	10000 U	
	C565	DIBENZOFURAN		
	C570	2,4-DINITROTOLUENE	10000 U	

#### KEYSTONE DC # --- 8

SAMPLE NUMBER: SPECOO18-02 INT-11C

SEMIVOLATILE ORGANICS ANALYSIS DATA SHEET, CONTINUED

**DATAFILE: 4U04243C05** 

 COMPOUND		DETECTION AMOUNT LIMIT FOUND (MILLIGRAMS / KG)	
C580 C585 C590 C595 C610 C615 C625 C630 C645 C645 C715 C725 C725 C725 C740 C740 C740 C740 C7775 C780 C785 C785 C790	DIETHYL PHTHALATE 4-CHLOROPHENYL PHENYL ETHER FLUORENE 4-NITROANILINE 4,6-DINITRO-2-METHYLPHENOL N-NITROSODIPHENYLAMINE 4-BROMOPHENYL PHENYL ETHER HEXACHLOROBENZENE PENTACHLOROPHENOL PHENANTHRENE ANTHRACENE DI-N-BUTYL PHTHALATE FLUORANTHENE PYRENE BUTYL BENZYL PHTHALATE 3,3'-DICHLOROBENZIDINE BENZO(A)ANTHRACENE BIS(2-ETHYLHEXYL)PHTHALATE CHRYSENE DI-N-OCTYL PHTHALATE BENZO(B)FLUORANTHENE BENZO(K)FLUORANTHENE BENZO(A)PYRENE INDENO(1,2,3-CD)PYRENE DIBENZO(A,H)ANTHRACENE BENZO(GHI)PERYLENE	10000 U 50000 U 50000 U 10000 U	410 J

U = UNDETECTED AT THE LISTED DETECTION LIMIT

J = COMPOUND IS PRESENT, BUT BELOW THE LISTED DETECTION LIMIT

SAMPLE: SPECOO18-02 INT-11C

ORGANICS ANALYSIS DATA SHEET - PAGE 5

LABORATORY NAME: KEYSTONE ENV. RESOURCES CASE NO . --

QC REPORT NO.: 05/18/92 ANALYST: JZ DATAFILE. 4U04243CO5

B. TENTATIVELY IDENTIFIED COMPOUNDS

CAS # SEMIVOLATILE COMPOUND NAMES SCAN# PURITY AMOUNT

MG/KG

NO NON-HSL COMPOUNDS FOUND > 10% OF NEAREST INT. STD

ESTIMATED VALUE - A 1:1 RESPONSE FACTOR IS ASSUMED